

UNIVERSITY OF GRONINGEN

BACHELOR THESIS

ASTRONOMY

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# Measuring the radio spectra of compact radio sources

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## Abstract

A study is performed into the angular extensions and differential number densities of gigahertz-peaked spectrum (GPS) compact radio sources (CRS) at low flux densities. An introduction is given into the topics of active galactic nuclei (AGN), CRSs and GPS, followed by a discussion of physical mechanisms abound in these sources. The selected sample of 45 CRSs is discussed and spectrally fitted using least squares and Markov-chain Monte Carlo algorithms, finding about half of the sources to indeed be GPS. Use is made of the K-z relation for radio sources, relating K-band magnitude to redshift, which is used to obtain redshifts for 14 sources without spectroscopic data, leading to a redshift completeness level in the sample of 66%. Number counts of redshift, peak flux density and peak frequency are made, showing no signs of bimodal behaviour. Obtained results are synthesised with data from Snellen et al. concerning angular size, which is shown to indeed decrease as expected at low flux densities and high frequencies. An unexpected offset is discovered in the small angular size-regime, which might hint at unusual behaviour of extremely young CRSs. Differential number counts of the number density per peak flux are performed, leading to the conclusion that a new population of unexplored, faint GPS sources may exist at low radio frequencies. Concluding remarks are provided as to how to proceed on this exploratory path, suggesting additional observations at differing frequencies, a redo of the Snellen et al. sample and exploration of the extremely high and low frequency ends of the known CRS range.

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# 1 Introduction

Before one delves into the subject of compact radio sources, it is imperative to gain a basic understanding of the physics behind astronomical radio sources in general, and the active galactic nuclei (AGN) that lie at their roots. To this end, this chapter will discuss the physical and observational backgrounds regarding cosmic radio sources, gigahertz-peaked spectra and previous research into this topic. At the end of this chapter a short summary of this thesis will be given.

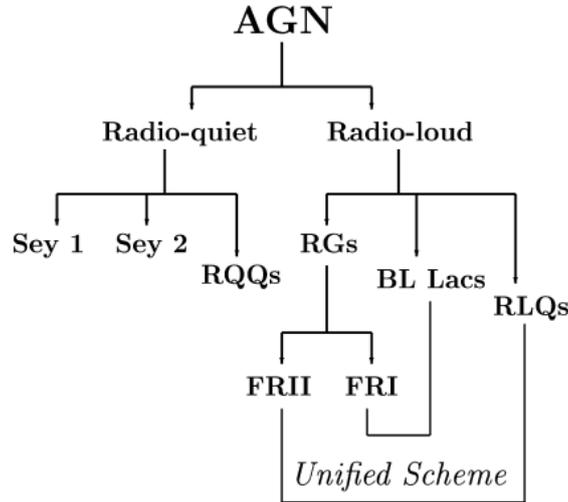


Figure 1: The classification scheme for AGN as presented by P. Kharb in her 2004 thesis. Sey stands for Seyfert galaxy, RQQs for radio-quiet quasars, RGs for radio galaxies, RLQs for radio loud quasars, BL Lacs for BL Lacertae objects and FR I and II for Fanaroff-Riley type I and II radio galaxies. Save for the last two types, none of the above objects will be discussed in much detail. The Unified Scheme refers to the hypothesis put forth by Barthel in 1989 that unifies radio galaxies with quasars [Kharb, 2004].

## 1.1 Radio sources & active galactic nuclei

The radio band of the electromagnetic spectrum is generally taken to range from  $\sim 100$  GHz down to the MHz regime or lower [Wilson et al., 2012]. It is an important scientific window into the cosmos, as the atmosphere is transparent to radio waves, thus enabling us to make radio observations from the Earth’s surface, in contrast to for instance infrared or X-rays. That radio waves are a relevant source of astronomical knowledge is in part because such a wide array of astrophysical processes emit them, or emit radiation that by its arrival to Earth has redshifted into the radio regime.

That radio waves had an extraterrestrial origin other than the Sun was first shown by Karl Jansky in 1931 [Jansky, 1933], who detected Sagittarius A\*, the galactic centre, and since then radio astronomy has taken a prominent place among the ranks of the astrophysical sciences. It has become one of the premier methods by which astronomers study young, distant and often luminous galaxies and their active galactic nuclei (termed AGN). These are nuclei of exceptional brightness (luminosities can exceed  $10^{46}$  erg/s) which can outshine the stellar light of their host galaxies. These sources are highly variable, on the order of days has been observed, and generally extend no further than several astronomical units (AU). They exhibit continuum emission over a large range of frequencies (low radio regime to X-ray), and their emission lines are strongly Doppler broadened, indicating rotational velocities of the gas in these AGN of up to  $10^4$  km/s, which orbit a central and highly active supermassive black hole (SMBH). This black hole powers the AGN by accretion of matter onto it, which by conservation of angular momentum is then released into perpendicular jets [Kharb, 2004].

There are multiple types of galaxies that house an AGN. Kharb presents a clear-cut scheme in her 2004 thesis which shows the dichotomy between these species (see Figure 1). As our thesis treats radio-loud AGN, and in particular radio galaxies, radio-quiet AGN will not be discussed in-depth here. The division between these two families of active galaxies is due to the difference in radio band output, with

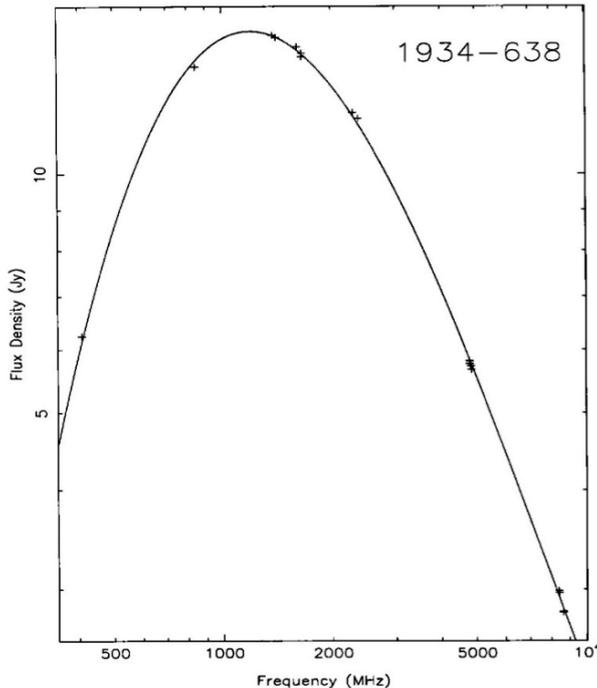


Figure 2: A quintessential GPS-shaped spectrum, belonging to PKS 1934–638, observed between 300 MHz and 8 GHz. This object is positioned at redshift 0.18, peaking just above 1 GHz. Taken from [Sadler, 2016].

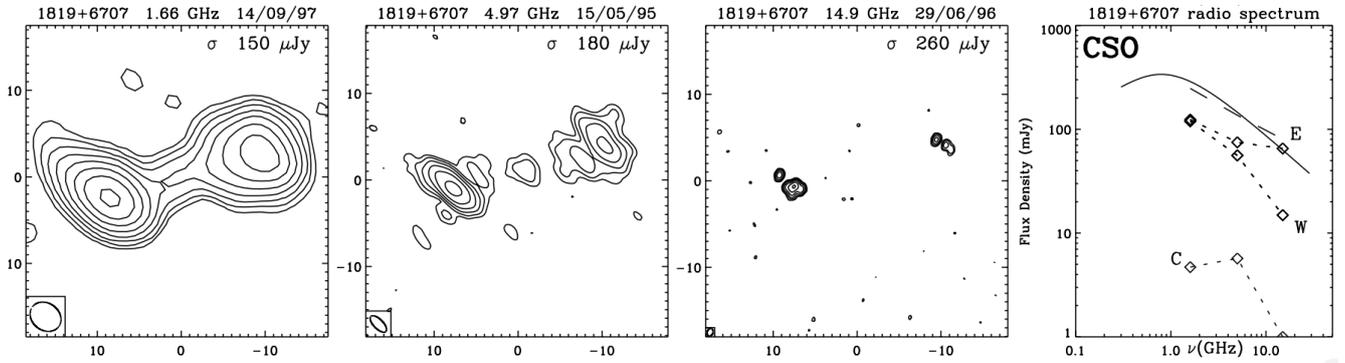
radio-loud galaxies, as the name implies, showing strong emission in radio. AGN are typified by the presence of a bright, star-like core with prominent emission lines, which are absent in normal galaxies. Radio-loud galaxies, however, have additional radio emission that can be an order of magnitude larger than the optical, and possess jets of highly energetic particles and ionized material which can extend up to hundreds of kiloparsecs perpendicular to the accretion disk around the central SMBH. This radio emission is commonly interpreted as synchrotron radiation originating from a non-thermal distribution of relativistic ( $v \sim c$ ) electrons [Kharb, 2004].

## 1.2 Evolution of compact radio sources

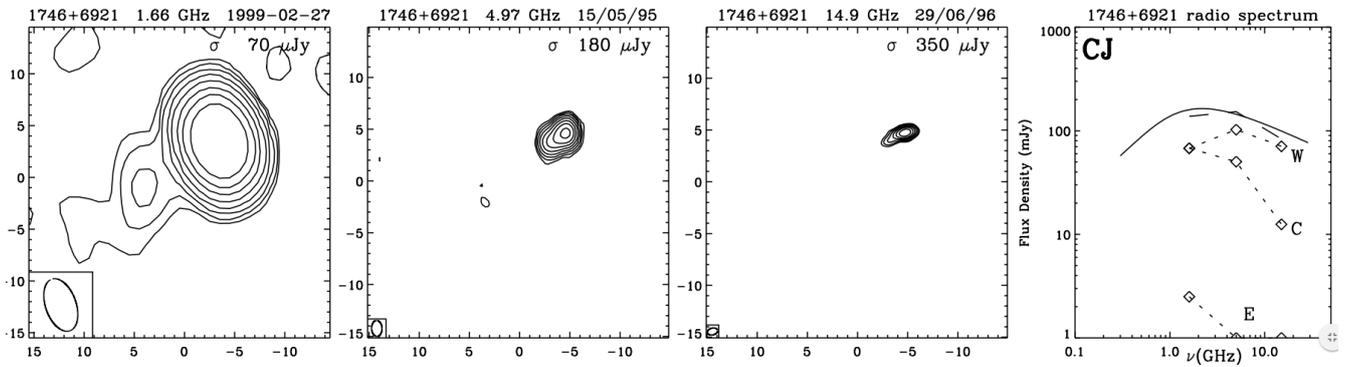
A specific subset of astronomical radio sources is made up by the so-called compact radio sources (CRS). These sources are often characterised by a small angular size (no larger than 1-2 arcsec), convex spectra that powerfully peak in the radio band, with a steep thick spectral index below the turnover peak and a thin spectral index of  $\alpha \sim 0.7$  above it [Orienti, 2016]. We will henceforth refer to sources as 'pure' if their spectra fit the above description fully, and as such have a form which is most similar to the theoretical spectrum of synchrotron emission by a homogeneous object showing signs of self-absorption at the low frequency-end (also known as a 'canonical' or 'classical' spectrum) [Mingaliev et al., 2013].

Compact radio sources can be divided into multiple categories, with the main divisions being spectral and morphological. Over the spectrum range, we find, from high to low frequency [Orienti, 2016]:

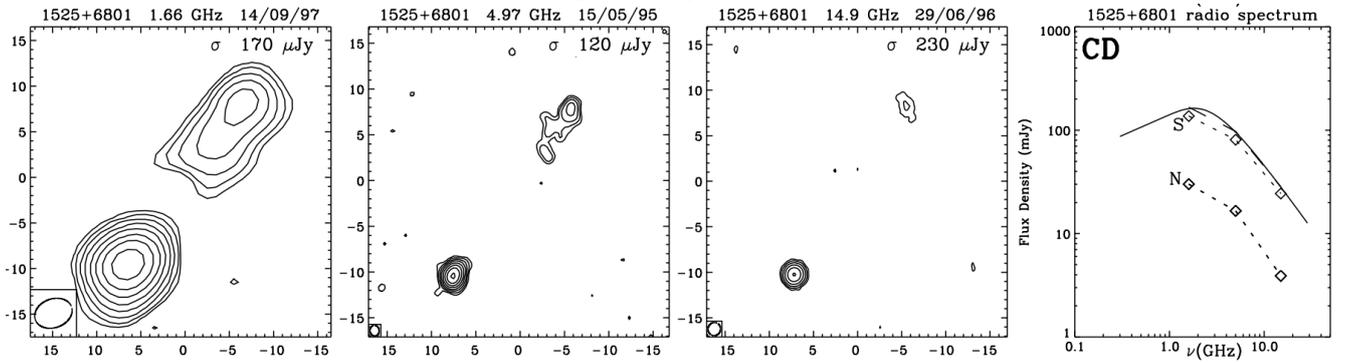
1. *High-frequency peakers (HFP)*: these are radio sources with their spectral turnover at the high end of the GHz-regime. They are considered a sub-population of the gigahertz-peaked spectra.
2. *Gigahertz-peaked spectra (GPS)*: these sources have a spectral turnover, as the name implies, in the GHz-regime, generally of the order of unity although it can vary from 0.5 to several GHz, as our sample will show. This thesis will look primarily at objects exhibiting these kind of spectra (see Figure 2 for a typical example).
3. *Compact steep spectra (CSS)*: this low-end type of radio spectra is typified by a spectral turnover in the high MHz-regime, typically of the order 100 MHz. As with HFP's and GPS, CSS show the same shapes of spectra, but at lower frequencies.



(a) A typical compact symmetric object (CSO).

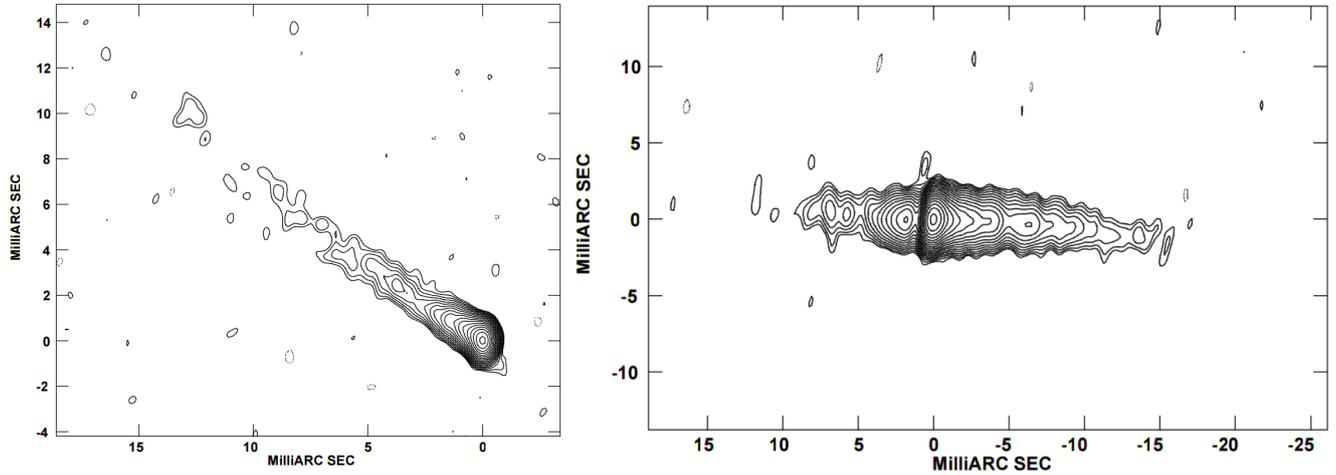


(b) A typical core-jet source (CJ).

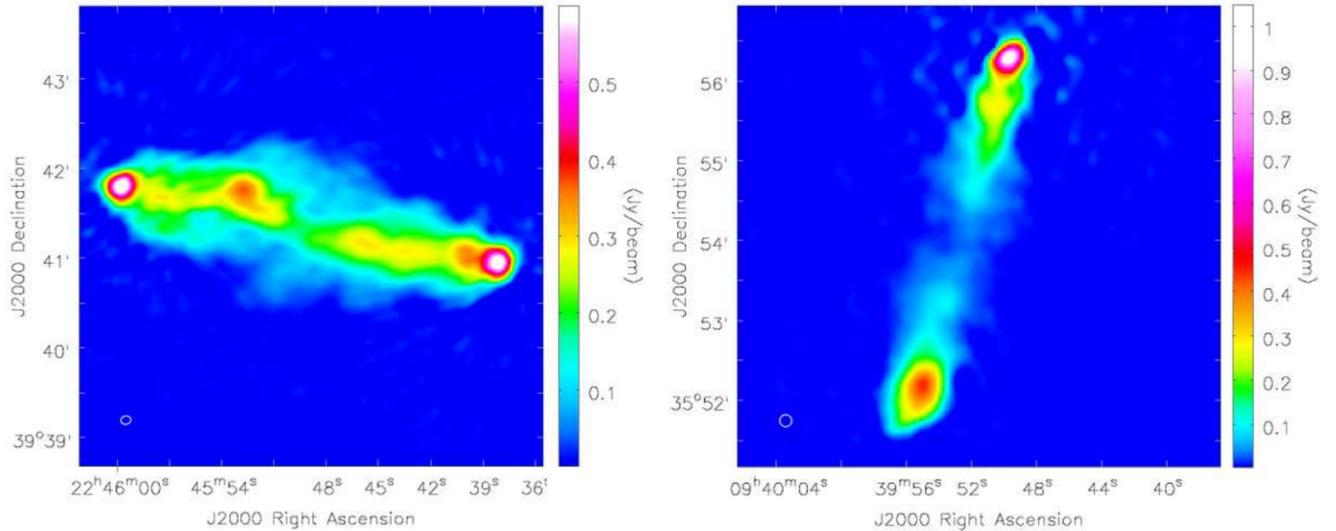


(c) A typical compact double (CD).

Figure 3: Examples of morphological classes of CRSS. Taken from [Snellen et al., 2000a].



(a) Two examples of FRI radio galaxies. Left: 3C 66B, imaged by combining 1.4 GHz VLA and 5 & 8 GHz VLBI observations, right: 3C 270, imaged by combining 5 GHz VLA and 8.4 & 8 GHz VLBI observations. Notice the lack of hotspots and the fuzzy appearance of the jets. Taken from [Kharb, 2004].



(b) Two examples of FR II radio galaxies. Left: 3C 452 at 138 MHz, right: 3C 223 at 147 MHz. Notice the prominent hotspots and well defined appearance of the jets. Taken from [Harwood et al., 2016].

Figure 4: Examples of the Fanaroff-Riley dichotomy in radio galaxies.

This division proves useful when one is interested in one particular stage of the evolutionary process of CRS's, as it has been implied that all three spectral classes are evolutionary stages a CRS passes through in the above order as it ages until it has grown enough to no longer be considered compact [Orienti, 2016]. In addition to this separation of types, there is also a morphological classification scheme (based on [Snellen et al., 2000b] and [Orienti, 2016]) - in particular order:

1. *Compact symmetric objects (CSO)*: these are objects with a milli-arcsecond scale symmetrical structure, often characterised by two-sided lobes (although in many cases one of the lobes is far brighter than the other), extending no larger than 1 kpc. They contain a flat spectrum component and have steep spectral slopes on either side (see Figure 3a).
2. *Medium-sized symmetric objects (MSO)*: if the source extends up to 10-15 kpc, they are termed medium-sized. Morphologically they look similar to CSOs.
3. *Core-jet sources (CJ)*: these sources are typified by a compact flat spectral component with a steeper component on only one side (see Figure 3b).
4. *Compact double (CD)*: sources showing two approximately equal components with similar spectra, but no indications of a central flat spectrum component (see Figure 3c).
5. *Complex sources (CX)*: sources possessing spectra that do not fall in any of the above categories.

Because the spectral and morphological divisions are so different in their criteria, selection of sources judging on the characteristics of one such category results in widely differing samples. However, Snellen points out that there is significant overlap between particularly GPS and CSOs: most of the GPS that are identified with galaxies have a mas-scale symmetric structure, and a large majority of CSOs possess a GHz-peaked spectrum [Snellen et al., 2000b].

A third division is the Fanaroff-Riley classification (FR), which divides radio galaxies into FRI and FRII types. The realization that radio galaxies came in two distinct species gave rise to this classification, which separates sources on the basis of the relation between core and jets: FRI sources show extended, diffuse tails which do not have clear termination points (see Figure 4a), while the jets of FRII sources are narrow, well collimated and terminate in hot spots (see Figure 4b) [Kharb, 2004]. FRII sources show bright lobes; FRI, conversely, are centre brightened [Wilson et al., 2012]. An additional category, FR0, was introduced to indicate a candidate radio source [Sadler, 2016].

Multiple hypotheses were put forth by O'Dea, Baum and Stanghellini in 1991 in one of the first papers on the topic of GPS ("What are the gigahertz peaked-spectrum radio sources?", [O'Dea et al., 1991]) about the nature of the sources that produced these spectra. They postulated that they could either signify that they were ordinary radio sources which were being frustrated by dense ISM, which limited the growth of the source's jets, or that they were young radio sources, and that GPS were the progenitors of type FRI/II radio galaxies, into which they would over time evolve [Orienti, 2016], [Snellen et al., 2000a]. The general consensus is in favour of the latter hypothesis, deduced from dust spectral ageing and observed expansion speeds, although the frustration theory cannot be ruled out.

It is suspected that the often sharp peak of all these compact sources is due to the high density of synchrotron emitting electrons in the source, which participate in synchrotron self-absorption. Another striking feature is their - cosmically speaking - extremely young age, sometimes less than 10 kyr, further adding to the suggestion that these objects are in fact young radio sources akin to Active Galactic Nuclei (AGN) which in turn evolve into larger sources [Snellen et al., 1998b]. GPS sources are estimated to comprise approximately 10% of the bright radio source population, with the CSS sources accounting for another 30% [Sadler, 2016].

While CSS and GPS sources were long considered the least variable radio sources in the night sky, but this hypothesis was strongly adjusted when later research found that especially HFP and GPS type CRS vary strongly over their lifetimes, only displaying the traditional convex steep spectrum during flaring events. Only  $\sim 30\%$  of the HFP/GPS galaxies maintain their convex spectra, while CSS radio galaxies are rarely, if ever variable. HFP, GPS and CSS quasars, however, do display large variability in both overall flux density as in their spectrum.

### 1.3 Previous research

Earlier research into the topic of GPS was carried out by primarily Cristopher O’Dea [O’Dea et al., 1991], [O’Dea, 1998] and Ignas Snellen, who performed research into the radio properties [Snellen et al., 1998b], infrared and optical spectrum [Snellen et al., 1998a], classification [Snellen et al., 2000b] and evolution [Snellen et al., 2000a] of these sources. Compact radio sources however were already known to the scientific community for more than thirty years [Allen et al., 1962], [Conway et al., 1963], [Blake, 1970], Blake being the first to suggest that CRS’s were young, and a portion of them had been catalogued by Peacock and Wall in 1981 and 1982 [Peacock and Wall, 1981] [Peacock and Wall, 1982]. The study of these objects began in earnest when Rudnick and Jones subjected a sample of CRS’s to a polarization analysis in 1982 and found that the degree of polarization showed little correlation with the observed wavelength [Rudnick and Jones, 1982]. The first sample of GPS (in that they had been selected with a peak flux around 1 GHz) was documented in 1983 as candidates for high redshift objects and radio sources with a symmetry at the mas scale [Gopal-Krishna et al., 1983] and a second sample was introduced two years later [Spoelstra et al., 1985].

The first real attempts at explaining the spectra of these sources were made by O’Dea, Baum and Stanghellini in 1991, who were the first to assume synchrotron self-absorption to be the cause of the GPS’s unusual shape, and used this inference to draw conclusions about them, further adding that in at least some sources free-free absorption too plays a role. They also found that these objects lack prominent jets, instead finding small, parsec scale steep-spectrum emission (dubbed ”micro-lobes”), and that the sources are tightly confined. Another feature was the nature of the GPS, which were found to comprise both galaxies and quasars, with the galaxies having a more symmetric mas scale structure and also often participating in mergers. Further conclusions were made concerning the formation (the young source versus frustration scenario which is mentioned in the previous section) and the redshift, which was found to be mainly high (half with a  $z$  larger than 3) [O’Dea et al., 1991].

Snellen et al., in their 1998 paper (”A New Sample of Faint Gigahertz Peaked Spectrum Radio Sources”), were the first to describe the observed spectra using a synchrotron self-absorbed function of the frequency. This succeeded in describing typical GPS in a sample ranging in peak frequency from 500 MHz to  $>15$  GHz, but failed to fit power law spectral distributions or combined spectra (different types of synchrotron spectra will be discussed in the next chapter). They also concluded that GPS sources have lifetimes of approximately 250 times shorter than a typical large scale radio source, and that source luminosity must decrease by a factor 10 during its transition from GPS- to radio source. Redshift distribution in the sample hinders a direct interpretation of the source counts, and it is concluded that more modelling of radio source evolution is required [Snellen et al., 1998b].

O’Dea makes several additional arguments for the evolutionary hypothesis in his 1998 paper, including an inverse relation between linear size and radiated radio power ( $P \propto l^{-0.5}$ ), and the fact that the relation between number density and linear size is consistent with what one would expect for young, evolving radio sources. He structures the arguments in favour of the young hypothesis into four parts: that their morphologies are similar to large-scale sources, that there seems not to be enough gas in their environments to justify the frustration hypothesis, that in the absence of cold gas these sources can indeed expand at velocities some  $\sim 10\%$  of the speed of light, and that neither GPS nor CSS sources show a diffusely emitting halo, which is what one would expect were they confined for their lifetimes. He goes on to conclude that, combining his own data with other literature results, confines can be placed on the age of these sources, which he estimates to be intermittent on timescales of  $10^4 - 10^5$  yr.

### 1.4 Motivation for & summary of this thesis

A critical question that should be asked before scientific research is carried out is, quite simply, why? Why is the to be performed research relevant, and why is it necessary? In the case of this thesis, the answer is similarly straightforward. The earliest stages of AGN are poorly understood, and the mechanism that leads to their triggering forms an enticing mystery. Although this thesis might not resolve this issue single-handedly, it remains important, critical even, that thorough examinations of CRSs are carried out, given that these sources possibly represent the youngest spawns of the cosmic AGN population. Analysis of these compact sources might point out major flaws in our theoretical framework regarding the formation of AGN, or instead provide convincing support for the prevailing theories. In the end, it is the author’s humble hope that this thesis may contribute, even if ever so slightly, to mankind’s

indomitable pursuit of knowledge.

This thesis will first discuss the dominant physical mechanisms in the sources we are studying: synchrotron radiation, free-free absorption and synchrotron self-absorption. The different spectra detected among our sources will then be explained and analysed. A short discussion of the selected sample will then be given, specifying locations and criteria, after which a brief introduction to the numerical methods used - least squares and Markov-chain Monte Carlo - to obtain the spectral fits will be given. This thesis will proceed by discussing each fit in short and providing the tables with the fitting parameters. The following chapter will first give an introduction to the K-z relation to obtain redshift from K-band magnitude, which will then be applied to find the redshifts of 14 members of the sample which do not have spectroscopic redshifts. Combining spectroscopic and K-z redshifts, number counts will be made per redshift. Results will then be synthesised to obtain values for luminosity distance, angular size differential number count per peak flux band. This thesis will conclude with a discussion of the obtained results, after which ideas will be opted as to how to continue in this scientific direction and some concluding remarks will be given. Throughout this thesis, a standard cosmological model is assumed, with  $\Omega_M = 0.286$ ,  $\Omega_V = 0.7139$  and  $H_0 = 67.80 \text{ km s}^{-1} \text{ Mpc}^{-1}$ , the most precise value from PLANCK data.

## 2 Radio source emission mechanisms

Radio-band radiation is emitted by a wide array of astrophysical processes, both thermal (blackbody spectrum) and non-thermal. This chapter will focus on the specific non-thermal mechanism of synchrotron emission and synchrotron self-absorption. In all formulas treated for the rest of this thesis, the following symbols denote the following quantities unless stated otherwise:

- $c$  : velocity of light
- $e$  : charge of the electron
- $m_e$  : mass of the electron
- $E$  : energy
- $v$  : velocity of the electron
- $a$  : acceleration of the electron
- $B$  : magnetic field
- $S$  : flux density
- $\nu$  : frequency
- $\gamma$  : Lorentz factor, given by  $\gamma = \frac{E}{m_e c^2} = \left(1 - \frac{v^2}{c^2}\right)^{-1/2}$

### 2.1 Synchrotron radiation

Synchrotron radiation is non-thermal radiation emitted by electrons in a magnetic field. Electrons gyrate around the magnetic field lines of the jets and accretion disk. At velocities of  $v < c$ , this is called cyclotron radiation, but if the particles are accelerated to relativistic velocities this is known as synchrotron radiation. According to the Larmor formula,

$$P = \frac{2e^2}{3c^3} \cdot \gamma^4 a'^2, \quad (1)$$

wherein  $P$  is the power of the electron and  $a'$  the acceleration of the particle in the comoving electron frame. Ergo: a particle that moves with some acceleration  $a'$  in its own comoving frame, radiates with some power  $P$  in the rest frame. Following from the Einstein-Planck equations of relativistic motion of an electron in a (homogeneous) magnetic field,

$$\frac{d}{dt}(\gamma m_e \mathbf{v}) = \frac{e}{c}(\mathbf{v} \times \mathbf{B}), \quad (2)$$

with  $\mathbf{v}$  the vector velocity and  $\mathbf{B}$  the vector magnetic field, we obtain the formula for the acceleration as a function of the velocity perpendicular to the field lines and the field's strength for a single electron,

$$a_{\perp} = \frac{eB}{\gamma m_e c} v_{\perp} \quad \text{with} \quad \omega_B = \frac{eB}{\gamma m_e c}. \quad (3)$$

Wherein  $a_{\perp}$  is the acceleration perpendicular to the magnetic field,  $v_{\perp}$  the velocity perpendicular to the direction of  $\mathbf{B}$  and  $\omega_B$  the non-relativistic gyration angular frequency. Therefore, over all emitted frequencies,

$$\langle P \rangle = \frac{2e^4 B^2 v_{\perp}^2}{3c^5 m_e^2} \gamma^2 = \frac{2e^4}{3c^9 m_e^4} \cdot B^2 E^2 v_{\perp}^2. \quad (4)$$

Wherein all symbols take their aforementioned meanings. So the radiated power of a single synchrotron electron is proportional to the squares of the magnetic field strength, the velocity and the energy of the particle. Hence the total power emitted by the source is dependent on its magnetic field and the velocity and energy of the electrons, which in turn are related to their frequency.

The frequency dependence of the radiated electron power is given by,

$$P(\nu) = \sqrt{3} \frac{e^3 B \sin(\alpha)}{m_e c^2} F\left(\frac{\nu}{\nu_c}\right). \quad (5)$$

Wherein  $\alpha$  is the pitch angle (the angle between the magnetic field and the velocity),  $F$  the dimensionless synchrotron spectral function,  $\nu$  the frequency of the radiation and  $\nu_c$  the critical frequency.  $F$  and  $\nu_c$  are respectively given by,

$$F\left(\frac{\nu}{\nu_c}\right) = \frac{\nu}{\nu_c} \int_{\nu/\nu_c}^{\infty} K_{5/3}(\eta) d\eta, \quad (6)$$

$$\nu_c = \frac{3\gamma^2 e B \sin(\alpha)}{4\pi m_e c^2} = \frac{3}{2} \gamma^2 \nu_G \sin \alpha \quad \text{with} \quad \nu_G = \frac{eB}{2\pi m_e c^2} \quad (7)$$

Here  $\eta$  is some arbitrary integration variable, and  $\nu_G$  is the relativistic gyration frequency. From empirical evidence - cosmic ray measurements - we know that the number density of electrons  $N(E)$  behaves like a power law,

$$N(E)dE = K E^{-\delta} dE, \quad (8)$$

in which  $K$  is some constant with units of energy to the power  $-\delta$  per volume,  $\delta$  some power law index and  $dE$  the energy range under investigation, while we know very little about the value of the pitch angle  $\alpha$ 's distribution. However, we also know that,

$$\epsilon(\nu)d\nu = \frac{dE}{dt} N(E)dE = \langle P \rangle N(E)dE \quad \text{hence} \quad \epsilon(\nu) = \langle P \rangle N(E) \frac{dE}{d\nu}, \quad (9)$$

where  $\epsilon$  is the volume emissivity and the other symbols take their aforementioned meanings. Substituting equations 8 and 4 into the above relation, and using the relation given by,

$$E = \gamma m_e c^2 \approx \left(\frac{\nu}{\nu_G}\right)^{1/2} \cdot m_e c^2 \quad \text{hence} \quad \frac{dE}{d\nu} \approx \frac{m_e c^2 \nu^{-1/2}}{2\nu_G^{1/2}}, \quad (10)$$

where  $\nu_G$  is again the relativistic gyration frequency, we obtain as a final result for  $\epsilon(\nu)$  for a power law energy distributed electron population,

$$\epsilon(\nu) = \frac{2e^4}{3c^9 m_e^4} \cdot B^2 E^2 v_{\perp}^2 K E^{-\delta} \frac{m_e c^2 \nu^{-1/2}}{2\nu_G^{1/2}}. \quad (11)$$

This is a messy relation which can be more transparently rewritten to the proportionality of frequency, particle energy and magnetic field strength, given by,

$$\epsilon(\nu) \propto B^2 E^{2-\delta} (\nu \cdot \nu_G)^{-1/2}, \quad (12)$$

which, because of the relations between energy and frequency (equation 10) and relativistic gyration frequency and magnetic field strength (equation 7), then becomes a proportionality of frequency and magnetic field strength alone given by,

$$\epsilon(\nu) \propto B^2 \left(\frac{\nu}{B}\right)^{1-\delta/2} (\nu \cdot B)^{-1/2} \propto B^{(1+\delta)/2} \cdot \nu^{(1-\delta)/2}. \quad (13)$$

If we then define a spectral index  $\alpha = (1-\delta)/2$  and realize that flux density is proportional to intensity, which is proportional to volume emissivity, i.e.  $S(\nu) \propto I(\nu) \propto \epsilon(\nu)$ , we obtain the proportionality which forms the cornerstone of this thesis, namely the relation between synchrotron frequency and flux density of a power law energy distribution for a synchrotron emitting electron population,

$$S \propto \nu^\alpha, \quad (14)$$

wherein  $\alpha$  is some spectral index, defined separately in different regions of the synchrotron spectrum [Wilson et al., 2012]. For a purely synchrotron emitting population, the spectrum follows such a power law distribution, leading to the remarkable conclusion that the spectrum of a power law energy distribution in an electron population is itself a power law.

Synchrotron radiation is also a beamed phenomenon. This is because the electrons, which gyrate around and experience a net movement in the direction of the magnetic field lines, emit in a cone towards the propagation direction [Rybicki and Lightman, 1986]. As such, synchrotron emission from an AGN, which possesses extremely strong magnetic fields, is stronger in the direction parallel to the magnetic field's orientation.

## 2.2 Synchrotron self-absorption

In regions where the synchrotron radiation intensity and density are high, the emitted photons interact with the emitting electrons via synchrotron self-absorption: they are reabsorbed by the electrons, which then reach a more excited energy state. This happens in CRS's and as such in GPS sources. This phenomenon is called synchrotron self-absorption (SSA) as the synchrotron emitting electrons absorb their own emission, and no external absorber is in play [Torniainen, 2008]. This behaviour is inversely proportional to the frequency, meaning that below some frequency  $\nu_{peak}$ , SSA dominates over synchrotron emission and the spectrum decreases sharply with a theoretical spectral index (called the thick spectral index) of 5/2. This can be proven using the concept of effective temperature of the electron plasma, the temperature of a theoretical black body source that would emit the same amount of radiation as the plasma,

$$T_e = \frac{E}{3k}. \quad (15)$$

Wherein  $T_e$  is the effective temperature and  $k$  is the Boltzmann constant. Using the relation between energy and frequency and magnetic field strength, this can be rewritten to,

$$T_e = \frac{\gamma m_e c^2}{3k} \propto \sqrt{\frac{\nu}{B}}. \quad (16)$$

Wherein all symbols take their usual meaning. The specific intensity, which is the intensity of a theoretical black body source that would emit the same amount of radiation as the plasma, is defined when the effective temperature equals the brightness temperature  $T_b$  (the temperature the electron plasma would have if it were a blackbody) of the plasma as,

$$I(\nu) = \frac{2kT_b}{\lambda^2}, \quad (17)$$

and as the wavelength  $\lambda$  is proportional to one over the frequency  $\nu$ , we obtain,

$$I(\nu) \propto \nu^2 \sqrt{\frac{\nu}{B}} = \nu^{5/2} B^{-1/2}. \quad (18)$$

When we then once more recall that  $S(\nu) \propto I(\nu)$ , we obtain our final result for the flux density of a synchrotron self absorbed source,

$$S(\nu) \propto \nu^{5/2}. \quad (19)$$

[Wilson et al., 2012] In thermal processes, this thick index has the value 2, known as the Rayleigh-Jeans value, and thus showing that SSA is not a thermal process [Rybicki and Lightman, 1986]. This value is in theory constant regardless of the electron population's spectral index above the turnover peak, although in practice such a steep value has not been observed in previous research<sup>1</sup> [Torniainen, 2008].

As the source ages, the highest energy electrons radiate away their energy first, leading to the overall flux of the source to drop at high frequencies. This leads to a high-frequency 'ageing cutoff' where the spectrum drops sharply. When the energy of the electron plasma drops, so does its frequency threshold for when SSA starts to dominate, so the peak shifts towards the lower end of the spectrum. This is one of the main pieces of evidence that supports the hypothesis that HFP's, GPS and CSS are, in that order, stages of the evolution of a young radio galaxy, as the peak frequency decreases with each type [Oriente, 2016]. Additionally, at very low frequencies, below the SSA region, a cutoff has been theorized because the spectra of individual electrons start to interfere with the synchrotron radiation. This has not yet been physically observed [Wilson et al., 2012].

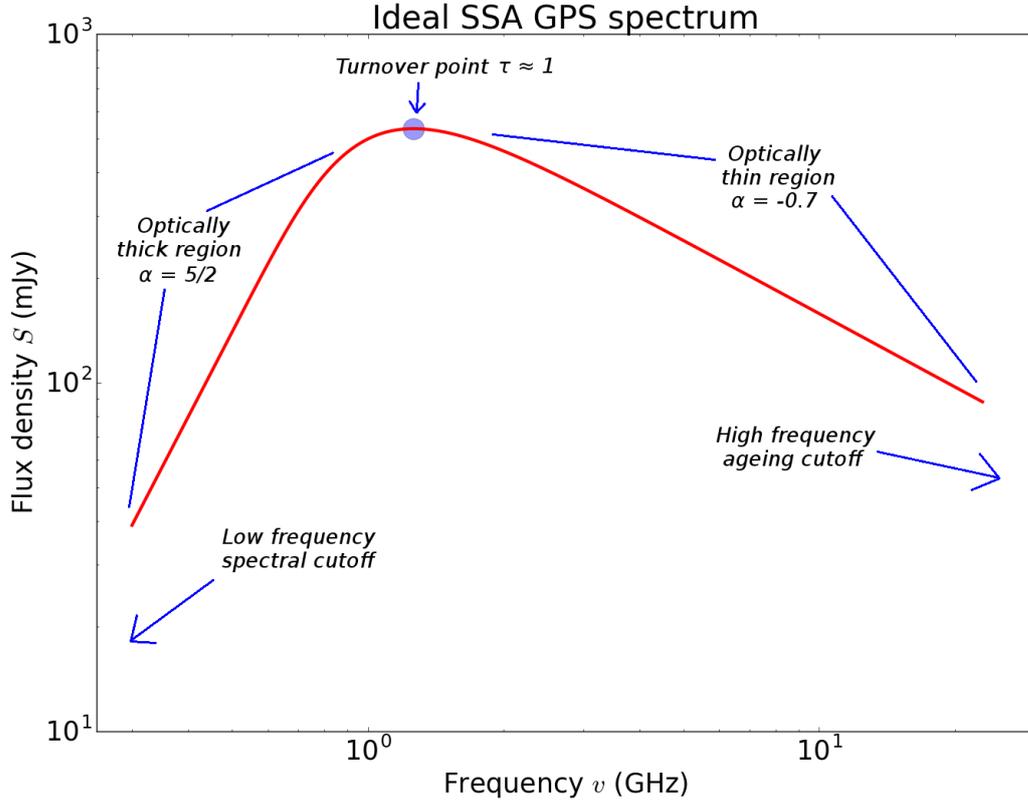
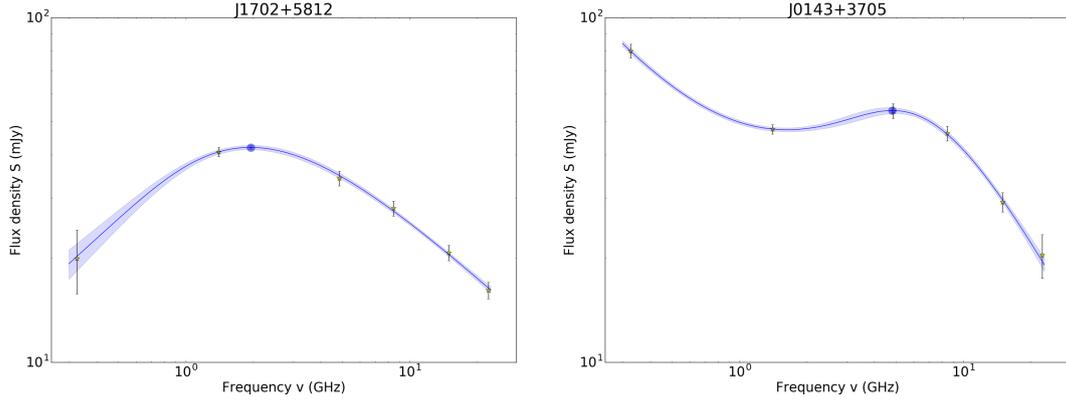


Figure 5: An annotated idealized synchrotron self absorbed GPS spectrum. Peak location at 1.3 GHz, peak flux 535 mJy, which exceeds the sources used in this thesis.

<sup>1</sup>In our sample, two sources (J0103+4322 and J0104+3840) possessed thick spectral indices of approximately 5/2. However, the author would like to point out that both spectra were notoriously hard to fit with a synchrotron peak, as the first of these two was almost a flat spectrum and the second one had a very poorly defined optically thick region, which was entirely described by just two data points, thus giving rise to large uncertainties.



(a) An archetypal GPS. Peak is located at 1.95 GHz, peak flux is 42 mJy. Spectral indices from Peak left to right: 0.6, -0.6 respectively.  
(b) A typical GPS+power law combined spectrum. Peak is located at 4.81 GHz, peak flux is 54 mJy. Indices: 0.5, -1.2, -0.8 respectively.

Figure 6: Examples of the two main types of peaked spectra.

## 2.3 Free-free absorption

Another process by which a strong turnover in the low GHz regime can be facilitated is free-free absorption (FFA). While free-free radiation (otherwise known as bremsstrahlung, after the German term for braking and radiation) is emitted when an electron in a plasma is slowed down and deflected by another charged particle, thereby losing energy in the form of a photon, during FFA the opposite occurs: the electron gains energy in its encounter with for instance an ion, absorbing a photon, causing it to jump to a higher energy state [Wilson et al., 2012]. The equation for FFA is given by,

$$S_{FFA}(\nu) = S_0 \cdot \nu^\alpha \cdot e^{\tau_f} \cdot \nu^{-2.1}. \quad (20)$$

Where  $S_{FFA}$  is the flux at at some frequency,  $S_0$  the scale flux density,  $\nu$  the frequency,  $\alpha$  the spectral index and  $\tau_f$  the optical depth at the peak. Suggestions have been made that in fact FFA by a thick gas- and dust disk around the AGN is responsible for the steep turnover in the lower frequency regions of a CRS spectrum, describing a “cold dense FFA plasma around the lobes of GPS sources” which “could be a cocoon which smothers expansion of jets and lobes” [Kameno et al., 2003], rather than SSA. If this were the case, then the degree of FFA, measured by the opacity  $\tau_\nu$  in equation 20, should vary with the viewing angle. Although Orienti comments that FFA cannot be completely discarded as an additional mechanism, because SSA spectra seem to fit the data accurately, more so than FFA, the scientific consensus leans towards SSA [Orienti, 2016], and as such the fits made in this thesis have been performed assuming SSA to be the dominant, if not only mechanism in play.

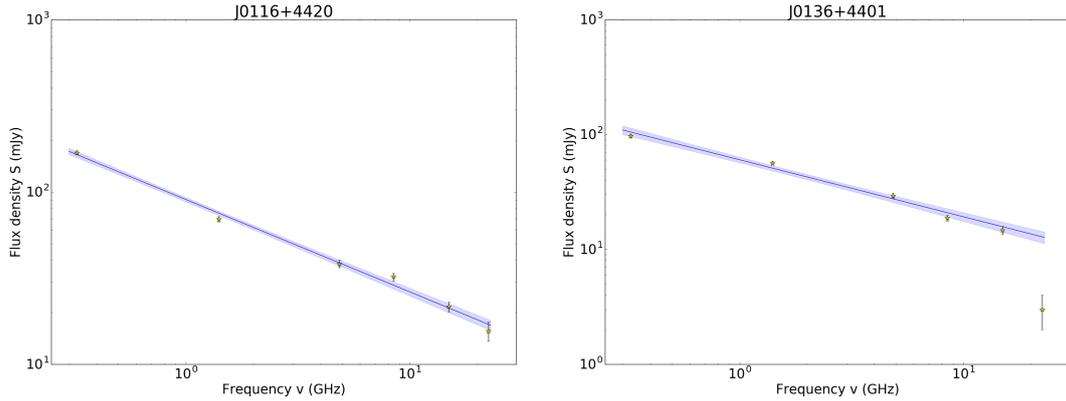
## 2.4 Synchrotron spectra

Although the quintessential SSA spectrum is given by the spectrum in Figure 5, multiple shapes of spectra appeared in the data, for which different fitting equations were used. These are discussed below.

### 2.4.1 Pure GPS spectra

Figure 6a shows a typical SSA GPS spectrum from our sample, with the shaded region denoting the one sigma regions around the fit. The fitting formula that we used in case of a typical GPS is given by,

$$S_{GPS}(\nu) = \frac{S_0}{1 - e^{-1}} \cdot \left(\frac{\nu}{\nu_0}\right)^k \cdot \left(1 - e^{-\left(\frac{\nu}{\nu_0}\right)^{l-k}}\right). \quad (21)$$



(a) A typical regular power law spectrum. Spectral index  $m$  is  $-0.54$ . (b) A typical aged power law spectrum. Spectral index  $m$  is  $-0.49$ .

Figure 7: Examples of the two main types of power law spectra.

This formula is taken from [Snellen et al., 1998b]. In optically thin regions (above the turnover peak), the spectral index, denoted by  $l$ , is negative, generally between  $-0.5$  and  $-1$ . This is because in these regions the radiation intensity and density are low enough that SSA by the plasma becomes negligible, and synchrotron emission can escape the source. Below the peak, the equation is dominated by the thick spectral index  $k$ , which as mentioned previously has a theoretical value of  $5/2$ , which represents a homogeneous, synchrotron self-absorbed radio source, but in practice generally varies between  $0.5$  and  $2$ . In consultation with Dr. McKean, this is explained by the spectrum being composed of multiple overlapping SSA components, some with a lower peak frequency and peak flux than the primary source, thus changing the spectral index  $k$  and making the increase more gradual in the low frequency regime (as is done in for instance [Kellermann, 1972]) [Orienti, 2016].  $S_0$  and  $\nu_0$  are scaling constants which influence the height and frequency of the peak, but have no further physical meaning.

Spectra with such a thick spectral index (i.e. when it is not exactly  $2.5$ ) are interpreted as multiple synchrotron emitting populations in different regions of the source, like nascent lobes, hotspots or jets. However, as only six data points were present across the observed band, a model with as many free parameters as a superposition of multiple synchrotron sources could not be fitted.

#### 2.4.2 Power law spectra

A second subtype of data sets was formed by the power law-fitted sets (Figure 7). The spectra of these sources were formed by power laws - or in some cases combinations of power laws - of varying spectral indices, generally in the same range as the thin spectral index of a synchrotron spectrum. The fit used for these kind of spectra is given by,

$$S_{pl}(\nu) = \sum_{i=1}^{n=1,2} S_{0,i} \cdot \left( \frac{\nu}{\nu_{0,i}} \right)^{m_i}, \quad (22)$$

in which  $m_i$  is the spectral index of the  $i$ -th power law population, and  $S_{0,i}$  and  $\nu_{0,i}$  the scaling coefficients of the  $i$ -th power law which determine the overall height and thus the size and intensity of that population - more strongly emitting populations have a larger height. The summation is used because linear combinations of power law spectra did occur,  $n$  being the number of power laws involved, but in only three occurrences where  $n = 2$ : in all other cases  $n$  was taken as unity. This is interpreted physically as an aged population of electrons which has passed through a synchrotron phase before but which has now radiated away most of its energy. Therefore, the turnover has shifted to such low frequencies that it is no longer within the observed range, and hence in the frequency window we use only a power law spectrum with  $n = 1$  is observed. SSA no longer plays an important role as synchrotron emission is no longer dense enough to enable it.

More than one type of spectrum necessitated a power law fit: The two main types are regular power laws (Figure 7a) and aged spectra (Figure 7b). These have a strong falloff at the high frequency end, with the data point at 22.49 GHz being offset negatively from the power law trend. Fitting an aged spectrum would be highly complicated and not required for the results to be obtained, so the decision was made to fit the remaining five points with a normal power law spectrum, effectively dropping the last point from the fitting process. These are physically interpreted as highly aged electron populations that are experiencing a high-frequency falloff as its most energetic electrons (and thus those with the highest energy) have radiated away.

### 2.4.3 GPS/Power law combined spectra

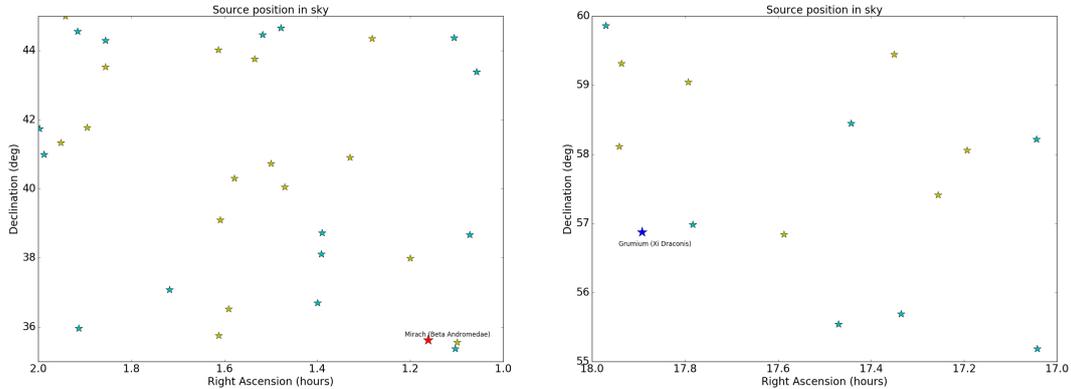
Spectra were also encountered wherein the first data point had a flux density offset from the optically thick part of the spectrum predicted by a simple SSA GPS fit (Figure 6b). These were fitted by assuming another, older population of electrons that follow a power law-distribution being superposed with the SSA spectrum, which gives rise to the combined formula,

$$S_{GPS+pl}(\nu) = S_{GPS}(\nu) + S_{pl}(\nu). \quad (23)$$

The physical interpretation of this combined population is a highly aged population of electrons in the source that have attained a power law distribution with a younger synchrotron component, implying that it has passed through a period of synchrotron emission before and that the present-day synchrotron component is not indicative of a young radio galaxy but an older, re-energized radio-loud source.

### 3 Fitting the radio spectra

The data sets obtained by McKean, Browne and Jackson showed clear synchrotron or power law behaviour but required fitting before any further conclusions could be drawn. This chapter discusses the methods used to obtain spectral fits to the used datasets and analyse the fit uncertainties. The entire procedure was carried out in the python programming language.



(a) The low field ( $1^h \leq \alpha \leq 2^h$ ,  $35^\circ \leq \delta \leq 45^\circ$ ). (b) The high field ( $17^h \leq \alpha \leq 18^h$ ,  $55^\circ \leq \delta \leq 60^\circ$ ). Mirach, a bright star, is provided for reference. Grumium, a bright star, is provided for reference.

Figure 8: The two observed fields. Cyan stars are sources who were proven to possess peaked spectra. Created using the script in Appendix B.4

#### 3.1 Sample screening and selection

The sample that has been used is drawn from data from observations done with the Very Large Array (VLA) at 4.86, 8.46, 14.94, 22.46 GHz. Measurements at 325 MHz and 1.4 GHz have been taken from the Westerbork Northern Sky Survey (WENSS) [Rengelink et al., 1997] and the NRAO VLA Sky Survey (NVSS) [Condon et al., 1998], respectively. The quasi-simultaneous VLA observations for each source were performed in 2000 and 2001 in C and D configuration while the WENSS and NVSS data were taken years before, resulting in in some cases large discrepancies between those data points and the VLA data and large uncertainty bars.

The sources were selected according to the following criteria:

- The spectral index in equation 14 between 1.4 (NVSS) and 4.85 (GB6) GHz satisfies  $\alpha \geq -0.5$ ;
- The flux density at 4.85 GHz according to the GB6 survey [Gregory et al., 1996] satisfies  $25 < S_{4.85} < 50$  mJy;
- The NVSS flux density at 1.4 GHz is measured by summing the radio emission within  $70''$  of the GB6 position;
- The flux density at 8.46 GHz as measured by CLASS [Myers et al., 2003] satisfies  $S_{8.46} \geq 16.7$  mJy.

The 45 sources of the complete sample are distributed between two fields: these regions are given by  $17^h \leq \alpha \leq 18^h$ ,  $55^\circ \leq \delta \leq 60^\circ$  and  $1^h \leq \alpha \leq 2^h$ ,  $35^\circ \leq \delta \leq 45^\circ$ , with  $\alpha$  and  $\delta$  the right ascension and declination respectively (see Figure 8). These two fields were chosen as to coincide with the fields used by Marlow et al. in his 2000 paper on spectroscopic redshifts [Marlow et al., 2000]: this becomes relevant in the determination of the K-z relation for this sample. The observation results are listed in Appendix A, with the VLA measurements in the second and the others in the first table. Sky areas have also been computed using the relation,

$$A_{sky} = (\alpha_{high} - \alpha_{low}) \cdot (\sin(\delta_{high}) - \sin(\delta_{low})). \quad (24)$$

## 3.2 Least squares fitting and Markov chain Monte-Carlo algorithms

Two numerical methods were used to compute the model parameters and estimate their respective uncertainties: least squares (LSq) and Markov chain Monte-Carlo (MCMC). These types of algorithm are shortly discussed in this section, but as most of the mathematical groundwork for these subroutines was already included in the modules from which the used functions were drawn, we will not go into great depths.

### 3.2.1 Least squares

Least squares fitting algorithms are a frequentist approach to model fitting and work by minimizing the sum of the squares of the differences between data and model over the uncertainty in the data (which can vary from point to point). Consider a model given by,

$$f(x) = a + bx. \quad (25)$$

Then the best fit to some data set  $y, x$  of size  $N$  is found by minimizing the quantity,

$$\chi^2 = \sum_{i=1}^N \frac{(y_i - f(x_i))^2}{\sigma_{y_i}^2}. \quad (26)$$

With  $\sigma_{y_i}$  the (Gaussian) uncertainty of the  $i$ -th data point. This quantity is called chi-squared, and the model is optimized for the data set when chi-squared is minimal. The advantage of this method is that numerically it is straightforward and reliable (especially if one is dealing with only six data points), but a pitfall rests in the fact that the minimum chi-squared obtained is in fact a false minimum: that is, somewhere else in the parameter space that is being sampled, there is a position with a smaller chi-squared [Hogg et al., 2010]. The covariances (and the variances, which essentially are simply a special case of the covariance where the two co-varying parameters are the same) are computed by,

$$Cov(X, Y) = \frac{\sum_{i=1}^N (x_i - \bar{x})(y_i - \bar{y})}{N}. \quad (27)$$

Wherein  $X$  and  $Y$  take the place of the parameters. This then constructs the covariance matrix, which in turn is used to compute the uncertainty in the actual fit. If more free parameters exist than data points, however, the system is said to be underdefined and no covariance estimates can be made (This stems from experience with the fitting algorithm, which will be discussed in the following sections).

### 3.2.2 Markov chain Monte-Carlo

Markov chain Monte-Carlo algorithms follow a more Bayesian approach to model fitting. It employs random sampling (the Monte-Carlo part) and in that random walking through the parameter space by a multitude of walkers (Markov chains). As taken from “Handbook of Markov Chain Monte Carlo”: suppose we have a program of the form,

```
Initialize x
repeat {
    Generate pseudorandom change to x
    Output x
}
```

This is the essence of MCMC: placing a walker in a phase space, performing a pseudo-random change<sup>2</sup> on that walker, and returning the new value. More exactly, when MCMC is used to fit a model to data, sampling is performed from an  $n$ -dimensional phase space of parameter values, where  $n$  is the number of parameters in the model (in the case of this thesis, this will vary from 2 to 6). To each point in this

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<sup>2</sup>Computer generated data is never truly random.

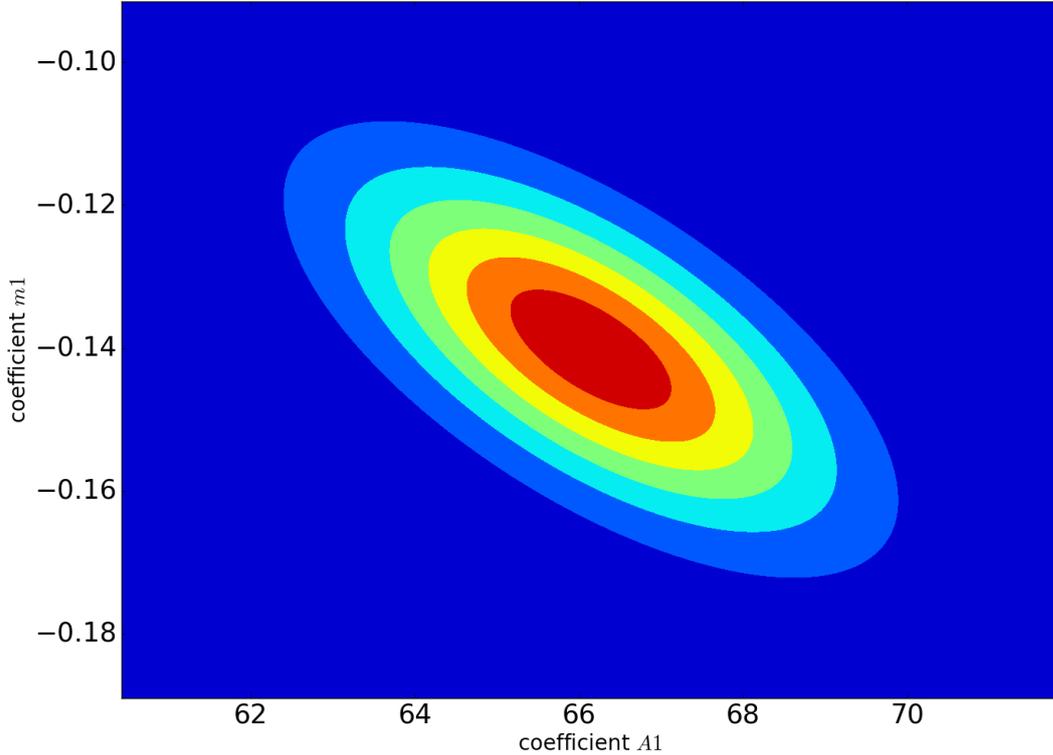


Figure 9: An example contour plot for the coefficients of a power law spectrum, in this case J0136+3905.

parameter space is assigned a probability value, and the optimal values for the model’s parameters are found by maximizing that probability: this maximization is performed by having the walkers return the probability value from the phase space after every random step. Is the probability value higher than the one found in the step before, then it is accepted as the new initial position for the next random iteration. If this is not the case, the walker falls back to its original starting position and attempts a different change. This continues until the walker finds a maximum from which it cannot escape. As we use a collective of multiple hundreds of walkers from which we take only the absolute highest value, the chances of collapsing into a false maximum are reduced. When one takes for instance the model in equation 25, the phase space of the parameters might look like the contour plot in Figure 9, where the region within the central isophote delineates the most probable values of the parameters within one sigma [Brooks et al., 2011].

In our code the *emcee* and *corner* python packages are utilized to perform the actual sampling. The *emcee* package takes a likelihood (the model) and prior (an indication of prior knowledge about the model, this can be used to shape the phase space to be biased towards certain values) function, distributes  $m$  walkers in  $n$  dimensions, then makes  $x$  iterations to their phase space position, eventually returning the minimum probability value found, which then provides the optimal parameters for the model. *corner* in turn uses the accrued walker values to create contour plots from which variances and covariances can be deduced by determining the full-width half-maximum (FWHM) values of the contours parallel to the parameter axes and the diagonals.

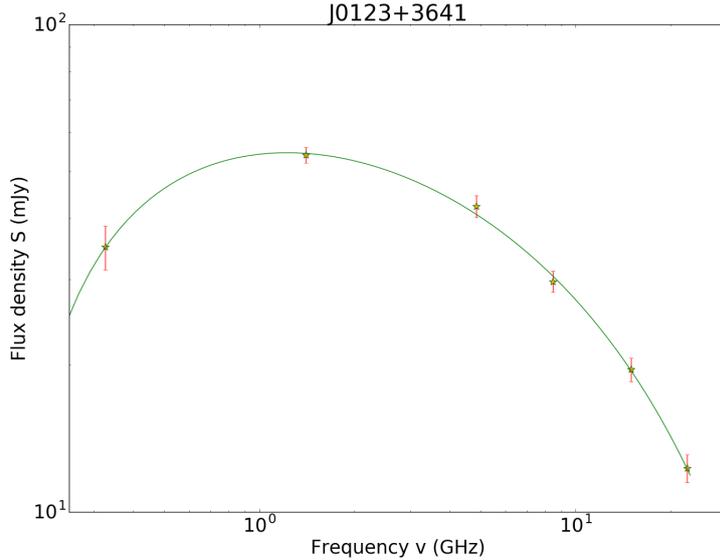


Figure 10: The polynomial fit to the spectrum of J0123+3641. This source was chosen to test the fitting algorithms as it was smooth, continuous and therefore easy to fit.

### 3.3 Methodology

#### 3.3.1 Polynomial fitting

A preliminary analysis of the spectra was performed by fitting them with a logarithmic polynomial (see Figure 10), given by the equation,

$$S(\nu) = \sum_{i=0}^{n=2,3} a_i \cdot (\log_{10}(\nu))^i. \quad (28)$$

This function has no physical significance and does not describe the synchrotron behaviour of the source’s electron population according to the model for our sources, but can be used as a preliminary analysis of the data and test the robustness of our code.

The function was described in python and subjected to an LSq algorithm from the *scipy.optimize* package to obtain the coefficients  $a_i$  for the polynomial by minimizing the sum of the squares of the uncertainties. Dependent on whether or not the data points showed signs of multiple turnovers, a quadratic ( $n = 2$ ) polynomial was fitted normally and a cubic ( $n = 3$ ) polynomial in case of two stationary points. As a polynomial proved substantially more intuitive and easy to fit than an SSA spectrum, this method was used to obtain a general impression of the sample (i.e. determine which spectra behaved like typical GPS, power law spectra or combinations of these) and find rough estimates for the peak frequency and peak flux. Extensive attention was not devoted to the uncertainties in either the coefficients or the peak, as this was not the final result and only served to give a crude impression of the spectral shape. This is also why an LSq algorithm was used instead of MCMC: LSq is substantially lighter and faster in terms of processing power and as the output served only as an estimate the risk of collapsing into a false minimum chi squared was accepted.

#### 3.3.2 SSA fitting

As a polynomial fit holds no physical meaning, the code was rewritten to fit an SSA spectrum to the data instead, as would physically be expected in the case of our sources. Although FFA fits to the data were also an option, the decision was made to focus on SSA. This spectrum is given by equation 21. In case a second turnover had been registered in the polynomial fit or was abundantly clear from the data first-hand, a GPS plus power law spectrum was fitted, given by equation 23. In case no turnover

had been registered, an ageing falloff was detected at the high frequency end or a negative turnover was detected, either one or multiple power laws were fitted, given by equation 22.

Flux density, flux uncertainty and observation frequency data are read from the table by the algorithm. These are then plotted to give the user an impression of the set, after which a user prompt enables picking one of the possible fitting procedures:

- *Power law spectrum*: This fit is selected if the spectrum does not show signs of curvature in log-log space. It includes all data points, which are then fitted with equation 22 for  $n = 1$ .
- *Aged power law spectrum*: This fit is used if the final data point shows a strong falloff with respect to the general trend through the other five data points. It excludes the final data point. The remaining points are then fitted with equation 22 for  $n = 1$ .
- *Double power law spectrum*: This fit can be selected if the spectrum shows signs of curvature, but concave instead of convex (which is what one would expect for a GPS). It includes all data points, which are then fitted with equation 22 for  $n = 2$ .
- *SSA GPS*: This is the type of fit used for stereotypical GPS spectra: convex, with a steep thick spectral index in the low frequency band and a more gently sloping thin spectral index in the higher frequency region. It includes all data points, which are then fitted with equation 21.
- *SSA+power law GPS*: This fit is applied if the spectrum shows a turnover peak, but curves back upwards in the lower frequency regions, i.e. it has two turnovers. It includes all data points, which are then fitted with equation 23.

The functions were then fitted using the `curve_fit` function from the `scipy.optimize` package, which applies an LSq algorithm to find the parameter values and estimates the covariances between them (see Appendix B.1). In case of either a power law or the SSA GPS fitting type, this succeeded. The function for a power law had to be rewritten for ease of coding to,

$$S(\nu) = \sum_{i=1}^2 A_i \cdot \nu^{m_i} \quad \text{with} \quad A_i = \frac{S_{0,i}}{\nu_{0,i}^{m_i}}. \quad (29)$$

As the degeneracies between  $S_{0,i}$  and  $\nu_{0,i}$  simply proved too great, leading to the fitting algorithm to give us results with massive uncertainties. The fitting of a function with simply an amplitude fudge factor  $A_{0,i}$  proved far more effective, and as the scaling coefficients had no real physical meaning to begin with this did not lead to non-physical results. The SSA+power law GPS is however deserving of some extra attention.

As only six data points were given per source, this frustrated efforts to determine the uncertainties of the parameters in a model with six degrees of freedom. LSq methods failed to account for this, so attempts were made to resolve this issue by instead performing the fitting procedure with an MCMC. To this end, the `emcee` and `corner` modules were acquired, which perform an MCMC algorithm and create covariance contour-plots, respectively. An MCMC fitting algorithm was coded (see Appendix B.2) using the values yielded by the LSq method as starting points for the walkers, assuming normal (Gaussian) prior distributions with small FWHM around these values in case of the spectral indices. 200 Walkers were used which were distributed randomly around the values estimated by the LSq method and then iterated over 1000 steps, returning values, uncertainties and a covariance matrix. The uncertainties in the parameters themselves and the covariances between them were determined by a function from the `numpy` package, but which utilizes the same equation as the LSq method (equation 27). To ascertain that the returned values were correct this was compared to the output of an algorithm that computes the uncertainties by determining the FWHM of the histogram of the relevant parameter as outputted by the MCMC. The `numpy.cov` function proved correct.

The peak and uncertainties in that peak were then found by using a bisection algorithm of a finite step size on the derivative of the fitting function - if the spectrum was fitted by a power law, then no peak was computed. The uncertainty in the peak flux was found by classical propagation of uncertainty (taking partial derivatives, multiplying by relevant covariance, square and sum for all parameters) while for the uncertainty in the peak frequency the step size of the bisection array was used. The spectrum

(with a shaded region around the best fit line indicating one sigma), contour plots and, in the MCMC's case, parameter histograms were then plotted and saved, as were the source name and index, parameter values for the best fits, their uncertainties and the peak coordinates, which were saved to a file.

### 3.3.3 MCMC uncertainty analysis (issues)

Although the MCMC was successful in finding the optimal values for the parameters to obtain the best fit line, the uncertainty analysis proved problematic. Uncertainties in the parameters for the GPS+power law fit were on average an order of magnitude in excess of what the LSq method returned, and when propagated through these uncertainties showed one sigma regions around the fit which exceeded the uncertainties in the data points themselves, something which was deemed non-physical. Suspicions were raised that this was due to high degeneracy between the spectral indices  $m$  (first power law, low frequency regime),  $k$  (second power law, mid regime) and  $l$  (third power law or exponential falloff, high end).  $m$  often proved poorly defined by only the two data points at 325 MHz and 1.4 GHz, thus prompting considerable uncertainties. Additionally, around 1.4 GHz the spectrum should transition from power law dominance, i.e. the  $m$ -regime, to dominance of the SSA component, and specifically the thick spectral region (the  $k$ -regime). This was assumed the culprit.

Attempts were thus made to combat this issue by strongly constraining  $m$ , first by taking a Gaussian prior with an extremely small FWHM, then by fixing it completely in the hopes that by allowing the program only five free parameters, this would push back uncertainties below the physical threshold of the data points' uncertainties. The formula used for fixing  $m$  is given by,

$$m = \frac{d[\log_{10}(S)]}{d[\log_{10}(\nu)]} = \frac{\log_{10}(0.5 \cdot S_{1.4GHz}) - \log_{10}(S_{0.325GHz})}{\log_{10}(1.4GHz) - \log_{10}(0.325GHz)}. \quad (30)$$

The factor 0.5 by which the flux at 1.4 GHz is multiplied stems from the fact that at in that frequency band the main contribution to the flux transitions from the power law component to the SSA component: therefore, at 1.4 GHz, half of the flux is contributed by the power law and half by the SSA spectrum. This fix proved ineffective. For reasons not discerned, uncertainties in the remaining five parameters remained an order of magnitude larger than outputted by the LSq method. When the MCMC was attempted on power law and SSA GPS spectra, with 2 and 4 free parameters respectively, success *was* achieved in obtaining uncertainties below the physically allowed values. Analysis of the problem showed that the mistake was not in the uncertainty propagation but stemmed directly from the MCMC output: the FWHM of the histograms confirmed the excessive uncertainties in the coefficients. This means that even after restraining one of the six parameters, degeneracies between the remaining five remain too high to provide a reliable result. As parameter uncertainties of the LSq method were of the same scale as for the MCMC in case of GPS and power law spectra, the LSq algorithm was used for these for time's sake and ease of computation. As in the GPS+power law case the MCMC returned non-physical uncertainties, another LSq algorithm was written to output the results for a GPS+power law fit with fixed  $m$ .

## 3.4 Results

The sample consisted of 45 sources. 44 Of these sources have flux density data at all six frequencies (0.325, 1.4, 4.86, 8.46, 14.94 and 22.46 GHz) and could thus be used. 23 spectra were peaked: 13 spectra were typical GPS and 10 behaved like GPS+power law combined spectra. The remaining 21 sources showed power law spectra, 5 of which showed signs of spectral ageing, 3 of which were double power laws<sup>3</sup> ( $n = 2$ ) while the remaining 13 behaved like regular power laws. Their coefficients and, in the peaked cases, peak coordinates are listed in the tables below the source-by-source discussion (Tables 1 to 6).

A full list of all the spectrum outputs is provided in Figures 11 and 12. A short discussion of each separate source will now be given in the same order as the given Figures. Some example corner plots of the covariance of the fit parameters will be provided to show what constitutes a good or a bad fit.

<sup>3</sup>The spectra fitted with double power laws were fitted as such as they clearly did not show a straight, power law like trend, but also lacked a clearly distinguishable peak, instead showing signs of a knee halfway through the frequency range. The fits here are somewhat arguable, as will become clear from the excessive uncertainties. However, all is not lost as these sources, being without peaks, are not used in the final results of this thesis.

1. *J0103+4322* shows signs of peaking in the high frequency regime, yet has an exceptionally flat background spectrum which can possibly be attributed to multiple overlapping SSA components. It took careful manipulation of the MCMC code to fit a GPS+power law spectrum to it as the thick spectral index  $k$  is extremely steep ( $\sim 2.5$ , the maximum physical value).
2. *J0104+3840* shows clear signs of a peak, but as the thick spectral index  $k$  was quite poorly defined by only the two non-VLA data points, the uncertainties there are rampant. This can be combated by tightly constraining  $k$ , but that would essentially be tampering with the results.
3. *J0105+3533* is an ordinary power law spectrum.
4. *J0106+3522* is an ordinary GPS.
5. *J0106+4422* shows GPS behaviour, but suffers similar issues as *J0104+3840* does: a poorly constrained  $k$ , in addition to an offset final data point, which might be indicative of a second, younger synchrotron component in the HFP regime.
6. *J0116+4420* is an ordinary power law spectrum.
7. *J0119+4054* is an ordinary power law spectrum.
8. *J0123+3842* is an ordinary GPS+power law spectrum.
9. *J0123+3806* is an ordinary GPS+power law spectrum.
10. *J0123+3641* is an ordinary GPS. This source was used to test the LSq and MCMC algorithms because of its typical shape.
11. *J0128+4003* is an ordinary aged power law spectrum. It was the first in our sample to be identified as such and thus sparked efforts to adapt the code to this kind of behaviour.
12. *J0128+4439* is an ordinary GPS.
13. *J0129+4044* is a double power law. Originally suspected a GPS+power law spectrum, after repeated attempts to find a peak failed the double power law solution was applied instead to much greater success.
14. *J0131+4428* shows GPS+power law like behaviour, but in this case the thin spectral index  $l$  proved to be under-defined. After adaptation of the code, this spectrum was successfully fitted, although the thick spectral index was outputted as  $\sim 2.5$ , an exceptionally high value.
15. *J0132+4345* is an ordinary power law.
16. *J0134+4018* is an ordinary aged power law spectrum.
17. *J0135+3631* is an ordinary aged power law spectrum.
18. *J0136+3905* is an ordinary power law.
19. *J0136+3545* is an ordinary aged power law spectrum.
20. *J0136+4401* is an ordinary aged power law spectrum.
21. *J0143+3705* is an ordinary GPS+power law spectrum.
22. *J0151+4332* is an ordinary power law, albeit with large scatter around the power law trend.
23. *J0151+4417* is an ordinary GPS+power law spectrum.
24. *J0103+4322* is a double power law. Attempts have been made to fit it with a GPS+power law spectrum but as  $l$  is barely defined this was unsuccessful, leading to a double power law with large uncertainties.

25. *J0154+3558* shows a GPS peak, but there are signs of an exceptional falloff in the high frequency regime. This could possibly be due to a highly aged background electron population overlaid onto the GPS.
26. *J0154+4433* is an ordinary GPS.
27. *J0156+4459* is an ordinary power law.
28. *J0157+4120* is an ordinary power law.
29. *J0159+4059* is an ordinary GPS+power law.
30. *J0159+4144* is a GPS which shows similar issues to J0154+3558.
31. *J1702+5511* is a GPS which shows similar issues to J0104+3840.
32. *J1702+5812* is an ordinary GPS.
33. *J1711+5803* is an aged power law, with signs of falloff at lower frequencies than usual.
34. *J1715+5724* is an ordinary power law.
35. *J1720+5541* is an ordinary GPS+power law.
36. *J1720+5926* is an ordinary power law.
37. *J1726+5826* is a GPS with small spectral indices. It suffers extreme degeneracies in its parameters (see Figure 17), presumably due to its small indices.
38. *J1728+5532* is a GPS with an unusually gently sloping  $k$ .
39. *J1735+5650* is an ordinary double power law.
40. *J1746+5659* is an ordinary GPS+power law.
41. *J1747+5902* shows signs of GPS+power law behaviour. However,  $l$  proved under-defined to such an extent that the spectral peak falls off the observed frequency band. Hence, although fitted with a GPS+power law model, it does not have a peak.
42. *J1756+5918* is a power law, but with a slightly offset 325 MHz measurement. This can be due to a small degree of SSA or FFA, but it was not enough for the LSq or MCMC of the SSA GPS model to pick it up.
43. *J1756+5806* is an ordinary power law.
44. *J1758+5951* is a GPS, but with an unusual amount of scatter in the data around the model's trend. As a result uncertainties are considerable.

In total some 10 sources have uncertain spectra, but these uncertainties do not affect later sections of the results. Therefore we take these spectra as final.

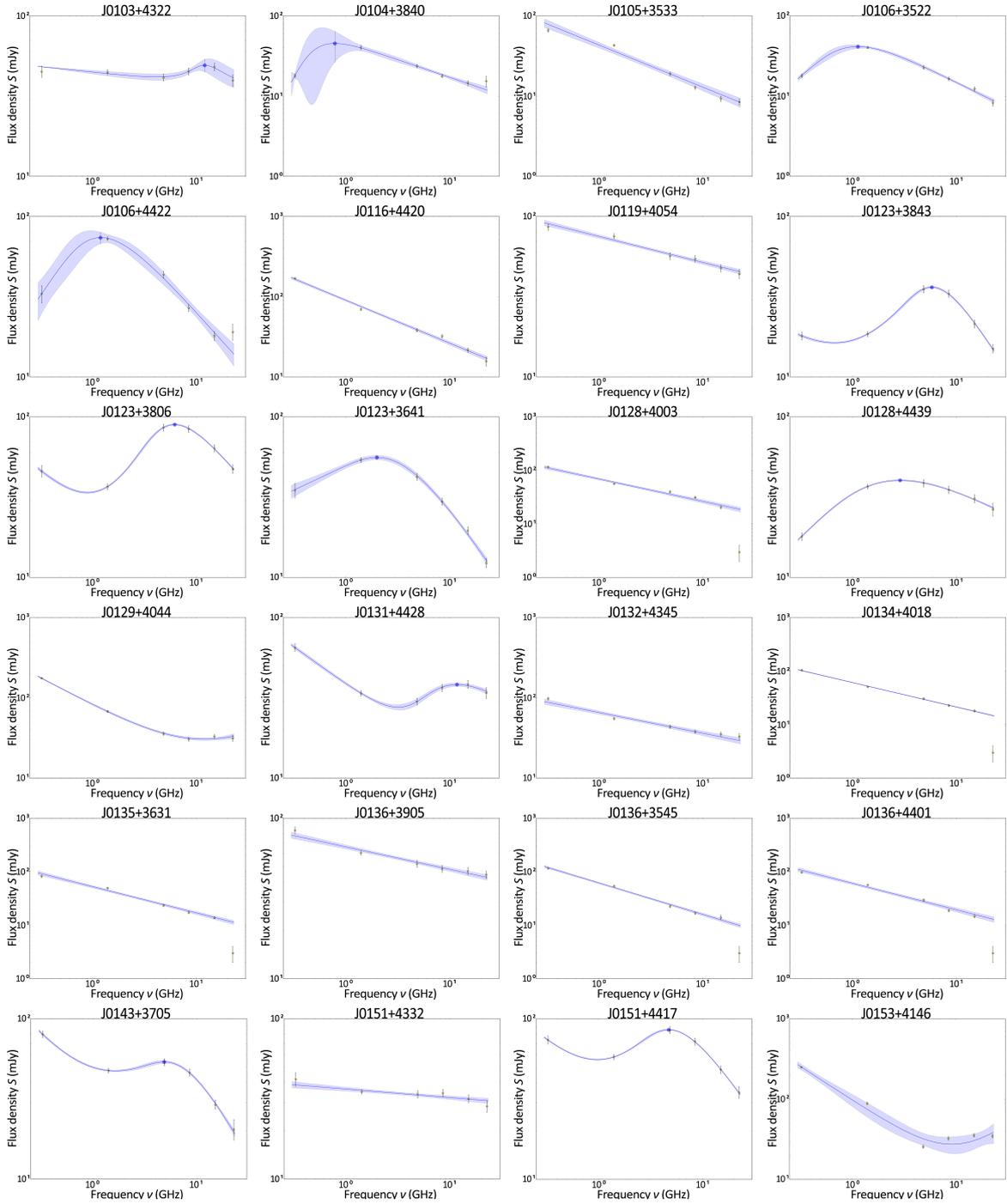


Figure 11: The first portion of the observed sources. List continues on the next page.

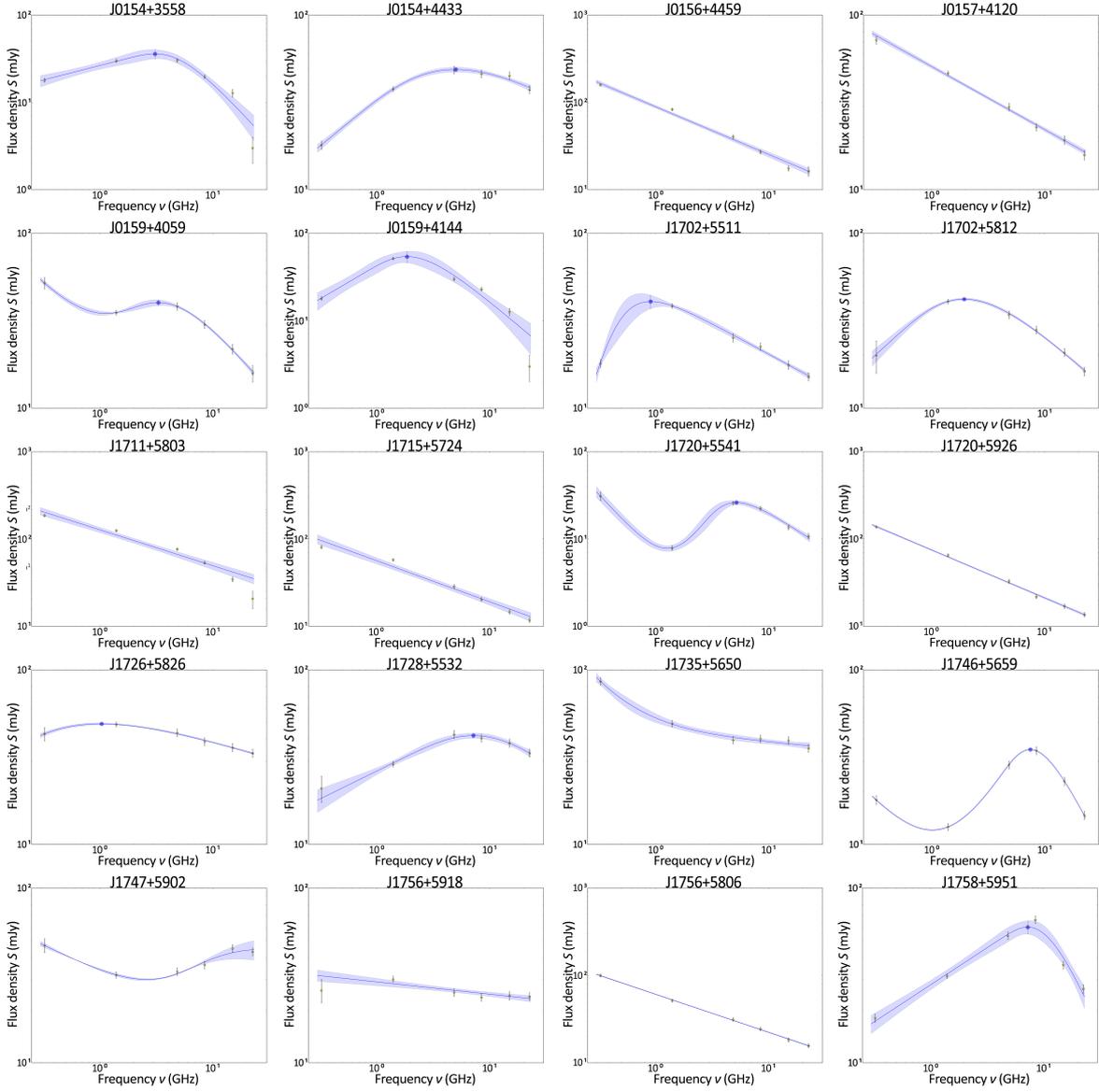


Figure 12: The second portion of the observed sources.

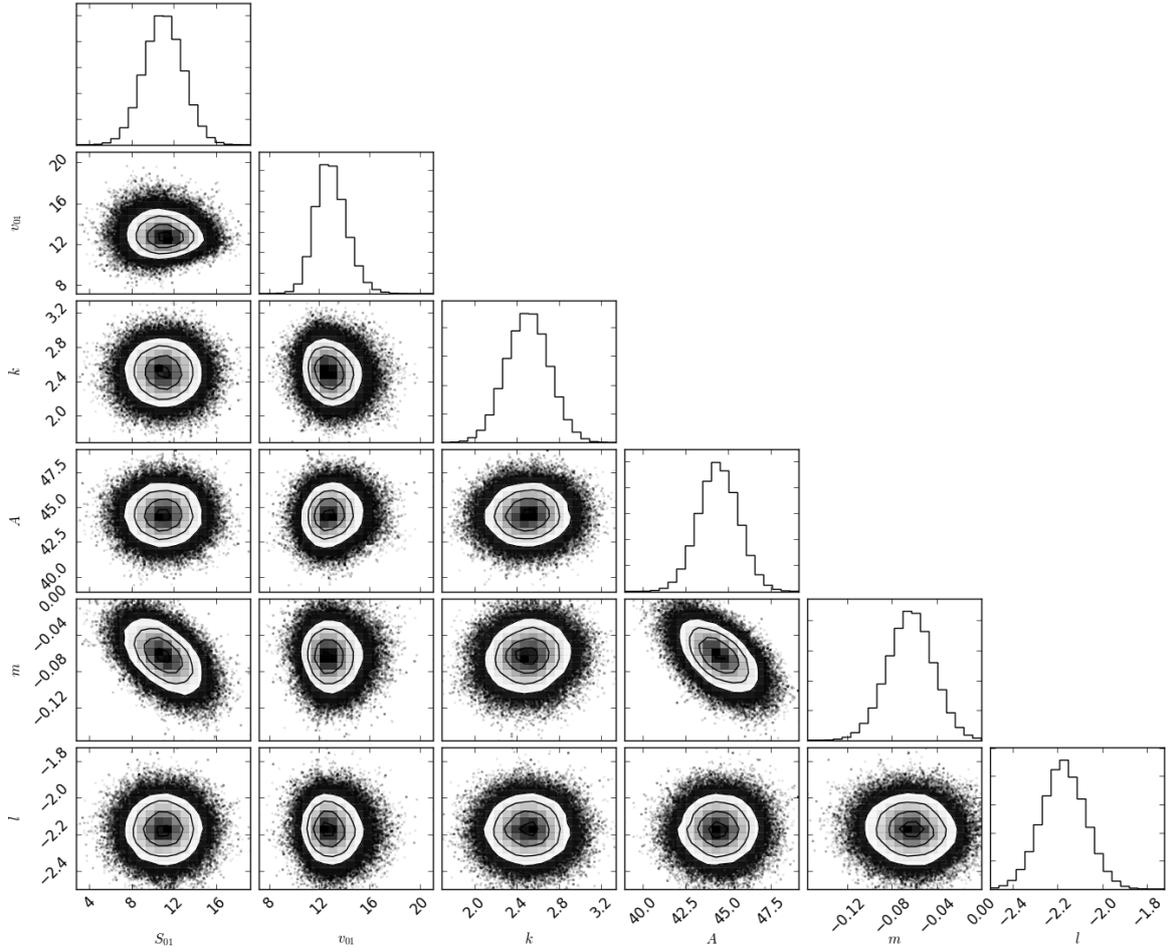


Figure 13: The contour plots of J0103+4322. Notice the relatively large uncertainties in  $k$  and  $l$ , contributing to the large final uncertainty in the fit. These contour plots were constructed from the MCMC algorithm (see Appendix B.2).

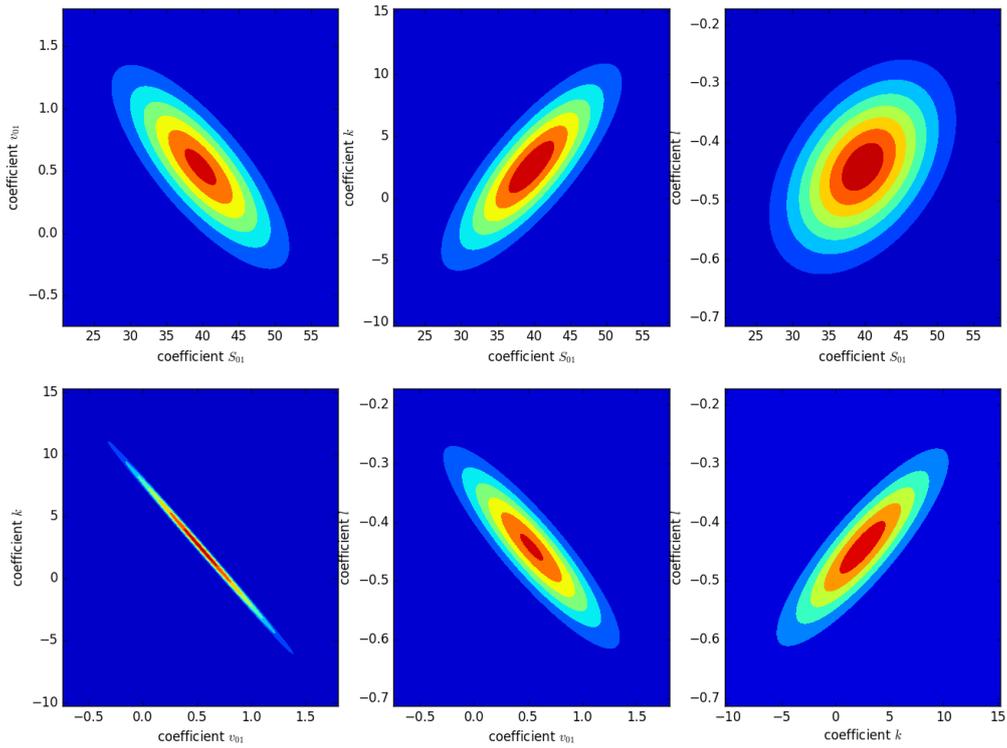


Figure 14: The contour plots of J0104+3840. Notice the major uncertainty in  $k$  and exceptionally high degeneracy between  $k$  and  $\nu_{01}$ , contributing to the large final uncertainty in the thick spectral region of the fit. These contour plots were constructed from the LSq algorithm (see Appendix B.1).

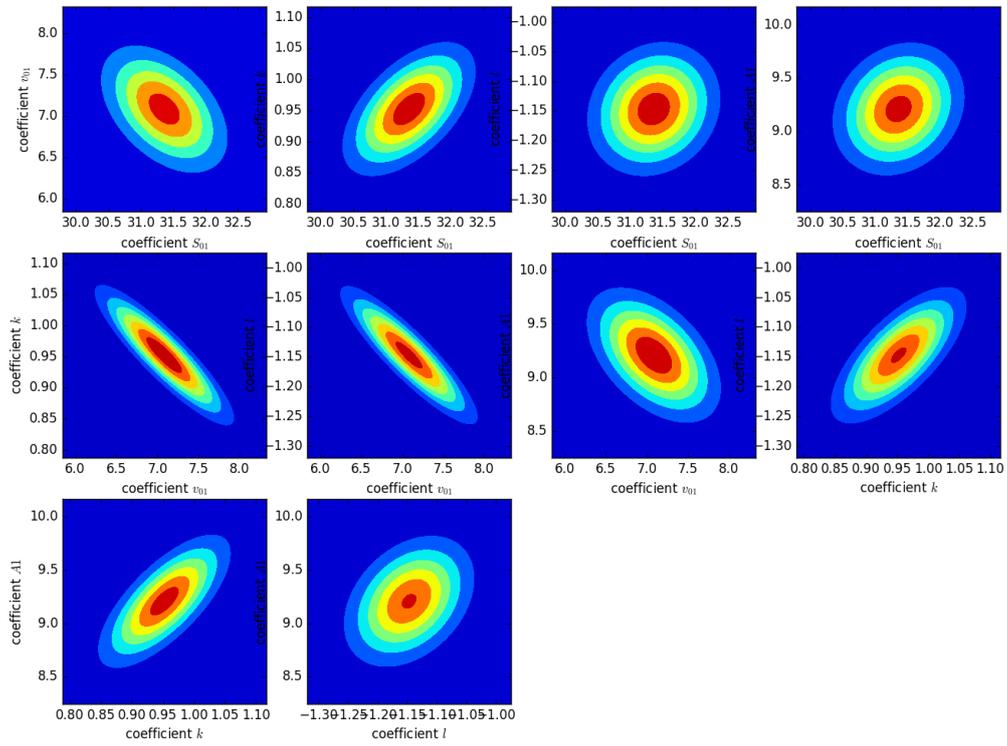


Figure 15: The contour plots of J0123+3842. These contour plots are an acceptable example for SSA GPS+power law fit contours: roughly circular shapes mean that there is only minor degeneracy, which reflects well in the uncertainties in the final fit. These contour plots were constructed from the LSq algorithm.

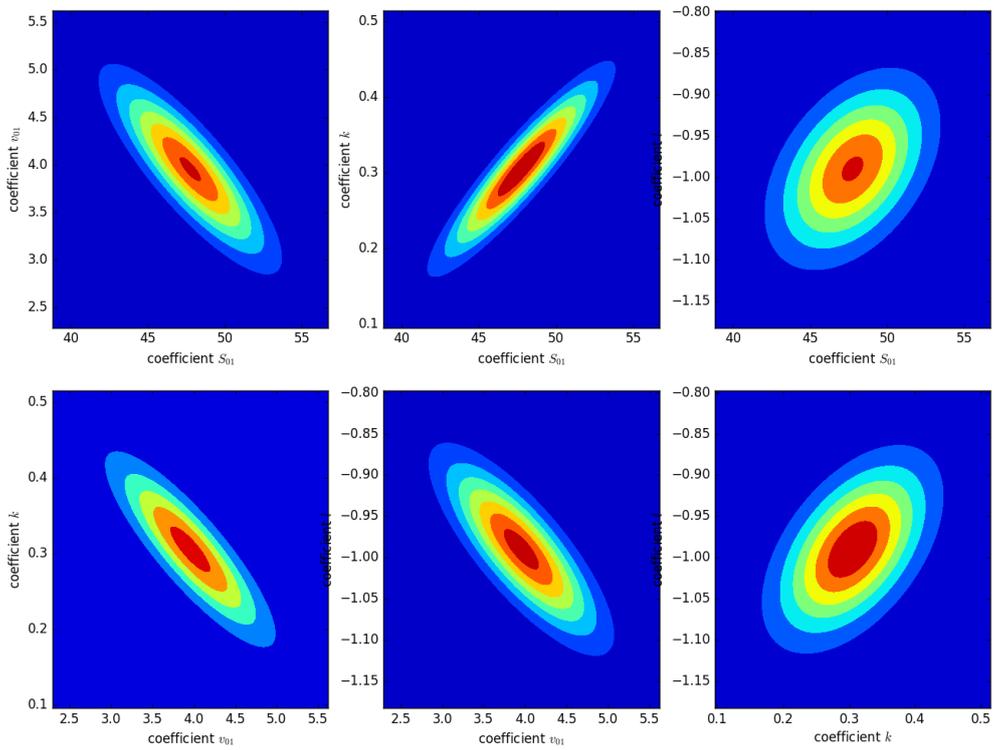


Figure 16: The contour plots of J0123+3641. Although the degeneracies are not minimal, these plots are an acceptable example for SSA GPS fits, as the magnitude of the uncertainties is small. These contour plots were constructed from the LSq algorithm.

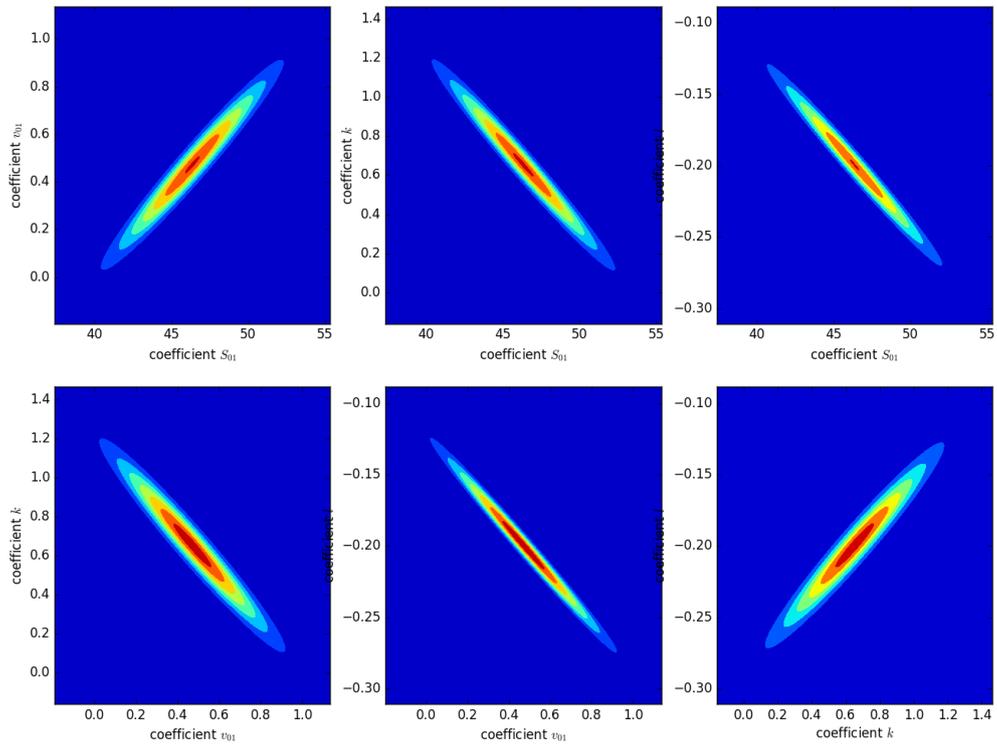


Figure 17: The contour plots of J1726+5826. As is clearly visible the degeneracies in these plots are significant, yet because of the small magnitude of the uncertainties it has little repercussions for the final fit. These contour plots were constructed from the LSq algorithm.

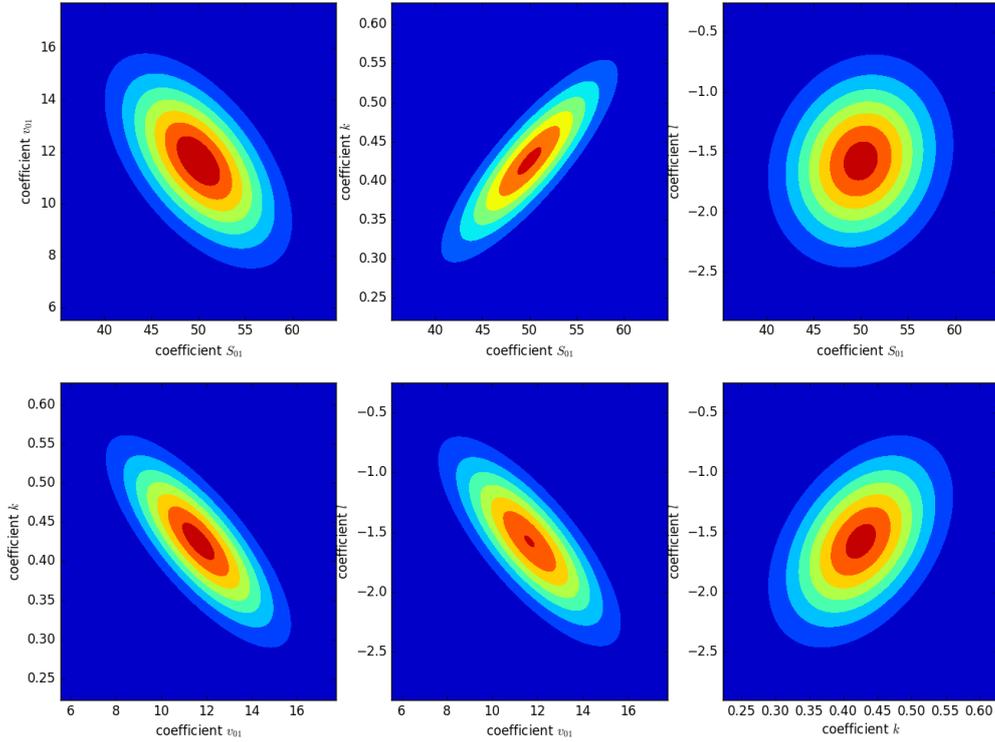


Figure 18: The contour plots of J1758+5951. While the spectral fit displays considerable uncertainties, degeneracies in these contours are smaller even than in the far more accurate cases of Figures 16 and 17, the uncertainties in the final fit are substantially larger, mainly due to sizeable magnitude of the uncertainties in the parameters themselves, leading us to conclude that degeneracies (and thus covariances) hold little to no direct relation to the quality of the fit. These contour plots were constructed from the LSq algorithm.

index	Source	type	$A$ (mJy GHz $^{-m}$ )	$m$
2	J0105+3533	regular	$43.4 \pm 3.5$	$-0.5 \pm 0.05$
5	J0116+4420	regular	$90.3 \pm 2.6$	$-0.5 \pm 0.02$
6	J0119+4054	regular	$75.0 \pm 1.9$	$-0.2 \pm 0.01$
10	J0128+4003	aged	$69.5 \pm 3.2$	$-0.4 \pm 0.03$
14	J0132+4345	regular	$65.5 \pm 3.2$	$-0.3 \pm 0.03$
15	J0134+4018	aged	$60.9 \pm 0.9$	$-0.5 \pm 0.01$
16	J0135+3631	aged	$52.3 \pm 2.9$	$-0.5 \pm 0.04$
17	J0136+3905	regular	$66.1 \pm 1.9$	$-0.1 \pm 0.02$
18	J0136+3545	aged	$61.2 \pm 2.1$	$-0.6 \pm 0.03$
19	J0136+4401	aged	$60.2 \pm 3.4$	$-0.5 \pm 0.04$
21	J0151+4332	regular	$36.4 \pm 1.0$	$-0.1 \pm 0.02$
26	J0156+4459	regular	$89.2 \pm 3.5$	$-0.5 \pm 0.03$
27	J0157+4120	regular	$50.8 \pm 1.2$	$-0.4 \pm 0.01$
32	J1711+5803	regular	$45.4 \pm 4.7$	$-0.6 \pm 0.06$
33	J1715+5724	regular	$56.3 \pm 4.6$	$-0.5 \pm 0.04$
35	J1720+5926	regular	$74.9 \pm 1.3$	$-0.5 \pm 0.01$
41	J1756+5918	regular	$29.0 \pm 1.4$	$-0.1 \pm 0.02$
42	J1756+5806	regular	$60.9 \pm 0.6$	$-0.4 \pm 0.01$

Table 1: The coefficients and their uncertainties for the 18 power law fits. The first column refers to the number index of that source within the python program used. The second column gives the source designation, the third specifies the subtype of power law, the fourth gives the power law amplitude and its uncertainty, the fifth gives the power law index and its uncertainty.

index	Source	$A_1$ (mJy GHz $^{-m}$ )	$m_1$	$A_2$ (mJy GHz $^{-m}$ )	$m_2$
12	J0129+4044	$79.6 \pm 22.2$	$-0.69 \pm 0.002$	$3.6 \pm 8.3$	$0.61 \pm 0.064$
38	J1735+5650	$9.8 \pm 170.5$	$-1.25 \pm 0.727$	$43.3 \pm 118.9$	$-0.05 \pm 0.007$
23	J0153+4146	$97.0 \pm 432.7$	$-0.85 \pm 0.037$	$1.3 \pm 24.1$	$1.01 \pm 1.379$

Table 2: The coefficients and their uncertainties for the 3 power law fits. The first column refers to the number index of that source within the python program used. The second column gives the source designation, the third specifies the subtype of power law, the fourth gives the power law amplitude and its uncertainty, the fifth gives the power law index and its uncertainty. The uncertainties here are exceptionally large - fortunately, spectral data for the double power laws are irrelevant for the final results of this thesis.

index	Source	$S_0$ (mJy)	$\nu_0$ (GHz)	$k$	$l$
1	J0104+3840	$39.7 \pm 6.4$	$0.53 \pm 0.42$	$2.50 \pm 4.26$	$-0.44 \pm 0.09$
3	J0106+3522	$41.2 \pm 1.7$	$1.03 \pm 0.31$	$1.12 \pm 0.32$	$-0.65 \pm 0.07$
4	J0106+4422	$73.7 \pm 6.4$	$1.25 \pm 0.77$	$0.94 \pm 0.52$	$-0.73 \pm 0.18$
9	J0123+3641	$47.8 \pm 3.0$	$3.96 \pm 0.56$	$0.31 \pm 0.07$	$-0.99 \pm 0.06$
11	J0128+4439	$40.1 \pm 0.5$	$2.41 \pm 0.30$	$0.63 \pm 0.04$	$-0.35 \pm 0.03$
24	J0154+3558	$29.3 \pm 5.0$	$5.43 \pm 1.37$	$0.33 \pm 0.12$	$-1.46 \pm 0.33$
25	J0154+4433	$48.8 \pm 1.6$	$4.96 \pm 1.24$	$0.53 \pm 0.06$	$-0.38 \pm 0.08$
29	J0159+4144	$52.0 \pm 7.8$	$2.37 \pm 0.93$	$0.76 \pm 0.23$	$-1.10 \pm 0.33$
30	J1702+5511	$35.5 \pm 1.1$	$0.56 \pm 0.17$	$2.06 \pm 1.11$	$-0.35 \pm 0.04$
31	J1702+5812	$41.6 \pm 0.8$	$2.37 \pm 0.33$	$0.59 \pm 0.09$	$-0.60 \pm 0.04$
36	J1726+5826	$46.4 \pm 3.0$	$0.47 \pm 0.22$	$0.65 \pm 0.27$	$-0.20 \pm 0.04$
37	J1728+5532	$38.6 \pm 3.4$	$14.08 \pm 5.86$	$0.32 \pm 0.09$	$-0.65 \pm 0.23$
43	J1758+5951	$50.0 \pm 4.9$	$11.65 \pm 2.04$	$0.43 \pm 0.07$	$-1.58 \pm 0.44$

Table 3: The coefficients and their uncertainties for the 13 SSA GPS fits. The first column refers to the number index of that source within the python program used. The second column gives the source designation, the third gives the scale flux and its uncertainty, the fourth gives the scale frequency and its uncertainty, the fifth gives the thick spectral index and its uncertainty, the sixth gives the thin spectral index and its uncertainty.

index	Source	$\nu_{max}$ (GHz)	$S_{max}$ (mJy)
1	J0104+3840	$0.79 \pm 0.02$	$45.3 \pm 18.6$
3	J0106+3522	$1.13 \pm 0.02$	$41.3 \pm 2.1$
4	J0106+4422	$1.20 \pm 0.02$	$73.7 \pm 6.1$
9	J0123+3641	$1.99 \pm 0.02$	$55.9 \pm 1.7$
11	J0128+4439	$2.88 \pm 0.02$	$40.3 \pm 0.4$
24	J0154+3558	$3.08 \pm 0.02$	$35.9 \pm 4.3$
25	J0154+4433	$5.01 \pm 0.02$	$48.8 \pm 1.6$
29	J0159+4144	$1.86 \pm 0.02$	$54.1 \pm 8.1$
30	J1702+5511	$0.90 \pm 0.02$	$40.7 \pm 3.6$
31	J1702+5812	$1.95 \pm 0.02$	$42.0 \pm 0.6$
36	J1726+5826	$1.04 \pm 0.02$	$49.1 \pm 0.6$
37	J1728+5532	$7.17 \pm 0.02$	$42.0 \pm 1.4$
43	J1758+5951	$7.26 \pm 0.02$	$59.8 \pm 5.2$

Table 4: The peaks and their uncertainties for the SSA GPS fits. The first column refers to the number index of that source within the python program used. The second column gives the source designation, the third gives the peak frequency and its uncertainty, the fourth gives the peak flux and its uncertainty. The peak frequency was taken as the step size of the bisection algorithm and is thus constant.

index	Source	$S_0$ (mJy)	$\nu_0$ (GHz)	$k$	$l$	$A$	$m$
0	J0103+4322	$11.5 \pm 1.8$	$12.75 \pm 1.24$	$2.52 \pm 0.20$	$-2.18 \pm 0.10$	$44.6 \pm 1.2$	-0.07
7	J0123+3843	$31.4 \pm 0.5$	$7.08 \pm 0.41$	$0.95 \pm 0.06$	$-1.15 \pm 0.06$	$9.2 \pm 0.3$	-0.46
8	J0123+3806	$82.8 \pm 1.1$	$6.09 \pm 0.35$	$1.29 \pm 0.06$	$-0.80 \pm 0.04$	$21.2 \pm 0.6$	-0.63
13	J0131+4428	$23.2 \pm 1.6$	$8.72 \pm 0.84$	$1.91 \pm 0.35$	-0.31	$38.3 \pm 1.0$	$-0.47 \pm 0.01$
20	J0143+3705	$41.6 \pm 1.7$	$8.39 \pm 0.85$	$0.48 \pm 0.06$	$-1.24 \pm 0.11$	$26.0 \pm 1.0$	-0.83
22	J0151+4417	$72.9 \pm 1.0$	$6.01 \pm 0.39$	$0.87 \pm 0.06$	$-0.98 \pm 0.04$	$31.5 \pm 0.7$	-0.64
28	J0159+4059	$32.2 \pm 1.4$	$4.52 \pm 1.36$	$0.81 \pm 0.28$	$-0.78 \pm 0.12$	$20.0 \pm 1.4$	-0.74
34	J1720+5541	$24.9 \pm 1.5$	$4.61 \pm 1.01$	$1.93 \pm 0.39$	$-0.85 \pm 0.17$	$6.3 \pm 0.9$	-1.41
39	J1746+5659	$32.4 \pm 0.2$	$8.80 \pm 0.10$	$1.10 \pm 0.01$	$-1.34 \pm 0.02$	$7.4 \pm 0.1$	-0.72
40	J1747+5902	$28.2 \pm 3.1$	$13.37 \pm 2.85$	$1.17 \pm 0.02$	$-0.12 \pm 0.01$	$32.1 \pm 1.9$	$-0.33 \pm 0.01$

Table 5: The coefficients and their uncertainties for the 10 SSA+power law GPS fits. The first column refers to the number index of that source within the python program used. The second column gives the source designation, the third gives the scale flux and its uncertainty, the fourth gives the scale frequency and its uncertainty, the fifth gives the thick spectral index and its uncertainty, the sixth gives the thin spectral index and its uncertainty, the seventh gives the power law amplitude and its uncertainty, the eighth gives the power law spectral index and its uncertainty. As  $m$  was fixed for most of the sources, it had no uncertainty. In case 13, instead  $l$  was fixed, and in case 40 neither was.

index	Source	$\nu_{max}$ (GHz)	$S_{max}$ (mJy)
0	J0103+4322	$12.07 \pm 0.01$	$49.2 \pm 4.4$
7	J0123+3843	$5.80 \pm 0.02$	$36.2 \pm 0.5$
8	J0123+3806	$6.20 \pm 0.02$	$89.6 \pm 1.0$
13	J0131+4428	$11.77 \pm 0.02$	$38.2 \pm 0.9$
20	J0143+3705	$4.81 \pm 0.02$	$53.7 \pm 1.1$
22	J0151+4417	$4.73 \pm 0.02$	$85.6 \pm 0.9$
28	J0159+4059	$3.29 \pm 0.02$	$40.1 \pm 1.5$
34	J1720+5541	$5.18 \pm 0.02$	$26.0 \pm 1.4$
39	J1746+5659	$7.49 \pm 0.02$	$35.0 \pm 0.2$
40	J1747+5902	N/A	N/A

Table 6: The peaks and their uncertainties for the SSA GPS fits. The first column refers to the number index of that source within the python program used. The second column gives the source designation, the third gives the peak frequency and its uncertainty, the fourth gives the peak flux and its uncertainty. The peak frequency was taken as the step size of the bisection algorithm, but as in the SSA+power law GPS case the bisection interval was shorter yet the number of steps constant, this varies slightly. In the case of index 40, this source behaved like an SSA+power law GPS but lacked a peak on the observed interval, hence the N/A's.

## 4 K-z relation

As not all GPS sources have well defined spectroscopic redshifts, another method has been devised which relates the K-band magnitude of a source to the redshift  $z$ , known as the K-z relation. This chapter discusses the K-band and subsequent redshift measurements of the sample, as well as the following calculations of luminosity distance and angular and linear size.

### 4.1 Introduction to the K-z relation

Light that travels over cosmic distances experiences a redshift  $z$  due to the cosmic expansion that causes the observed frequency of the emitted light to change via the relation,

$$z = \frac{\nu_{emit}}{\nu_{obs}} - 1 \approx \frac{v_{rad}}{c}. \quad (31)$$

Wherein  $\nu_{emit}$  is the frequency of the emitted light,  $\nu_{obs}$  the frequency of the observed light and  $v_{rad}$  the radial velocity of the source. This means that light from objects with a positive redshift and thus a positive radial velocity (i.e. away from us) is shifted towards lower frequencies.

The near-infrared K-band ( $\sim 2.2 \mu\text{m}$ ) magnitude of a radio galaxy is dominated by starlight from older stars and therefore associated with evolved ellipticals, although emission lines that have been redshifted into the K-band window also contribute to the band flux. The strength of these lines is intrinsically related to the underlying quasar continuum and hence the radio emission [Jarvis et al., 2001]. In our sample however, this contribution is negligible as the spectra show no lines.

Lilly and Longair found in 1984 that there was a correlation between the K-band magnitude and the redshift of radio galaxies. They found that the dispersion around this trend in the K-direction was approximately constant, Gaussian and quite tight out to redshift  $z \sim 2$ . The relation they found was not entirely in line with the expectations and attempts were made to explain this by assuming a changing number of red giant stars [Lilly and Longair, 1984]. Over time the dispersion along the K-band was narrowed down, first by Jarvis et al. to around 0.59, then by Cruz et al. to an improved significance of  $0.593 \pm 0.02 \text{ mag}$  [Jarvis et al., 2001], [Cruz et al., 2007], the value used in this thesis as well. Willott et al. in 2003 found the optimal fit through the data to follow the logarithmic quadratic polynomial function,

$$K(z) = 17.37 + 4.53 \cdot \log_{10}(z) - 0.31 \cdot \log_{10}^2(z) \quad (32)$$

Where  $K(z)$  is the magnitude in the K-band. This is the function we use in our study as well.

### 4.2 Sample selection

In order to obtain as complete a sample in terms of redshifts as possible, for those sources in the original sample of which a spectroscopic redshift was not known [Marlow et al., 2000], infrared measurements were taken using the United Kingdom Infrared Telescope (UKIRT) by McKean (2003). A limitation in the fact that only sources with a right ascension between  $01^h$  and  $02^h$  could be observed led to a final completeness of the sample (in terms of redshift by spectroscopy and K-band magnitude, which leads to K-z relation redshift) for 30/45 sources, or redshift completeness 66.6%. K-band measurements within 5 arcseconds are provided in Table 7.

### 4.3 Methodology

Using the *normal* function from the *numpy* module that draws random numbers from a Gaussian distribution given some mean and dispersion, five million data points were generated in  $K, \log_{10}(z)$ -space. This was done by first computing the mean in the K-magnitude by generating a wholly random  $\log_{10}(z)$ -value on the interval  $-1 < \log_{10}(z) < \log_{10}(5)$  using the *numpy.random* function, plugging this into the K-z relation and using the output as the mean for the Gaussian random number generator, which then returned a K-band magnitude. This created five million data points with a random  $z$  and a Gaussian dispersion around the K-z relation of value  $\sigma_K = 0.593$ , as required (see Figure 19 for the K-z relation

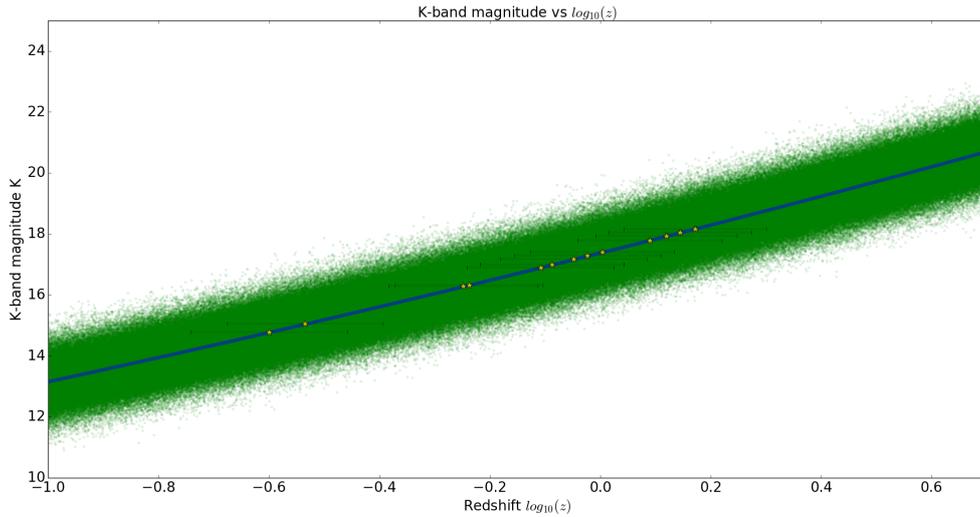


Figure 19: The K-z relation in  $\log_{10}(z)$  space. Yellow stars represent the points of which K-band data is known, the blue line is the idealized K-z relation according to equation 32 and the green points are the 5 million normally distributed, simulated data points.

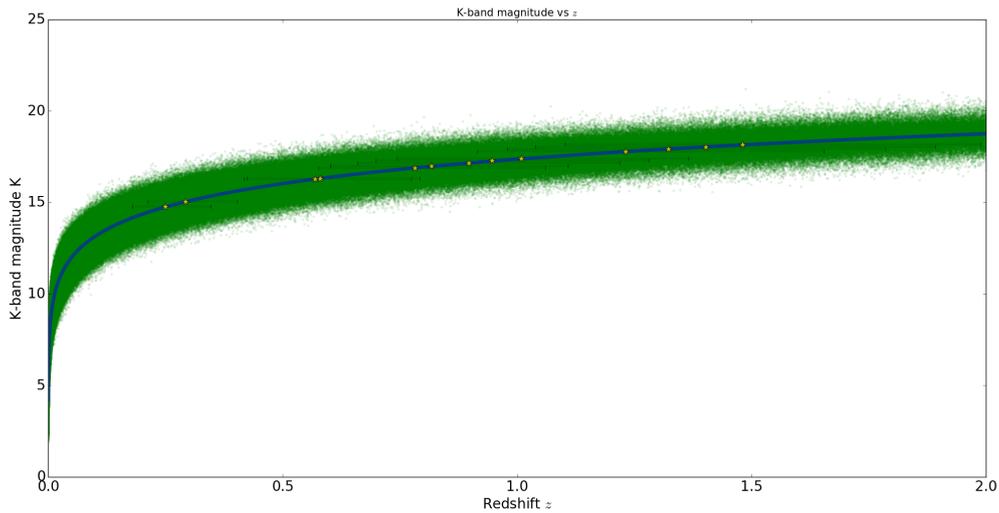


Figure 20: The K-z relation in linear  $z$  space.

in log-space and Figure 20 in linear space).

A sampler was then created which took a K-band magnitude and collected all of the random data points within a small interval around this K-magnitude (sample size  $\pm 0.05$  mag). Upon plotting the  $\log_{10}(z)$ -values within this interval in a histogram, this yielded a set of Gaussian distributions (see Figure 21), as one would expect from the near linear trend visible in Figure 19. When translated to linear  $z$  values this reproduced a log-normal distribution as expected (see Figure 22).

The mean ( $\mu_y$ ) and dispersion ( $\sigma_y$ ) of a Gaussian distribution in  $\log_{10}(x) = y$  are given by,

$$\mu_y = \frac{\sum_{i=1}^N y_i}{N}, \quad (33)$$

$$\sigma_y = \sqrt{\frac{\sum_{i=1}^N (y_i - \mu_y)^2}{N}}. \quad (34)$$

In which  $N$  is the length of the data set and  $y_i$  the  $i$ -th value in the data set. Alternatively, the corresponding statistics in the subsequent log-normal distribution in the variable  $x$  itself, namely the median and the lower and upper 68% confidence interval limits, are given by,

$$\mu_x = 10^{\mu_y}, \quad (35)$$

$$\sigma_{x,low} = 10^{\mu_y - \sigma_y} - \mu_x, \quad (36)$$

$$\sigma_{x,upp} = 10^{\mu_y + \sigma_y} - \mu_x. \quad (37)$$

Where  $\sigma_{x,low}$  and  $\sigma_{x,upp}$  denote the upper and lower uncertainties in  $x$ . They can also be computed directly from the log-normal distribution but this is substantially more convoluted so instead the data in  $K, \log_{10}(z)$ -space was used to compute the above statistics in the given sequence. The results were then outputted to a table and plotted, and the resulting K-band magnitude/redshift values were overlaid onto the K-z relation plot (see Figures 19 and 20, sources are indicated by yellow stars) (see Appendix B.3).

Source	spectrum	$K$ (mag)	$\log_{10}(z)$	redshift $z$
J0103+4322	GPS+PL	14.78±0.03	-0.6±0.14	0.25 <sup>0.10</sup> <sub>0.07</sub>
J0106+3522	GPS	16.99±0.04	-0.087±0.13	0.82 <sup>0.29</sup> <sub>0.21</sub>
J0106+4422	GPS	17.29±0.05	-0.02±0.13	0.95 <sup>0.34</sup> <sub>0.25</sub>
J0116+4420	PL	16.33±0.03	-0.24±0.13	0.58 <sup>0.21</sup> <sub>0.15</sub>
J0119+4054	PL	16.29±0.03	-0.25±0.14	0.57 <sup>0.21</sup> <sub>0.15</sub>
J0123+3843	GPS+PL	16.99±0.05	-0.087±0.13	0.82 <sup>0.29</sup> <sub>0.21</sub>
J0123+3641	GPS	18.17±0.08	0.17±0.13	1.48 <sup>0.51</sup> <sub>0.38</sub>
J0129+4044	DPL	17.94±0.08	0.12±0.13	1.32 <sup>0.45</sup> <sub>0.34</sub>
J0136+4401	APL	17.79±0.06	0.09±0.13	1.23 <sup>0.43</sup> <sub>0.32</sub>
J0143+3705	GPS+PL	17.17±0.04	-0.048±0.13	0.90 <sup>0.32</sup> <sub>0.23</sub>
J0153+4146	DPL	16.90±0.04	-0.11±0.13	0.78 <sup>0.28</sup> <sub>0.21</sub>
J0154+3558	GPS	17.40±0.05	0.004±0.13	1.01 <sup>0.35</sup> <sub>0.26</sub>
J0154+4433	GPS	15.05±0.03	-0.53±0.14	0.29 <sup>0.11</sup> <sub>0.08</sub>
J0159+4144	GPS	18.05±0.08	0.15±0.13	1.40 <sup>0.48</sup> <sub>0.36</sub>
J0123+3806	GPS+PL			1.656±0.004
J0128+4003	APL			3.525±0.003
J0128+4439	GPS			0.228±0.001
J0131+4428	GPS+PL			1.123±0.001
J0132+4345	PL			1.812±0.001
J0136+3545	APL			1.871±0.001
J0151+4332	PL			2.192±0.002
J0151+4417	GPS+PL			1.976±0.003
J0156+4459	PL			0.214±0.001
J0157+4120	PL			0.0811±0.0001
J1711+5803	PL			0.1465±0.0001
J1715+5724	PL			0.0273±0.0001
J1720+5926	PL			0.5878±0.0003
J1728+5532	GPS			1.404±0.002
J1747+5902	GPS+PL			0.981±0.001
J1756+5806	PL			0.192±0.002

Table 7: The results for the K-z relation fitting of the sources for which K-band magnitude data is known. As  $z$  is log-normally distributed, it has non-Gaussian upper and lower uncertainty boundaries. Sources in the lower section of the table have been determined spectroscopically (16 by spec., 14 by K-z) by Marlow et al. (2000). The second column indicates the fit type: GPS is a normal SSA GPS spectrum, GPS+PL is SSA GPS+power law, PL is regular power law, APL is aged power law, DPL is double power law.

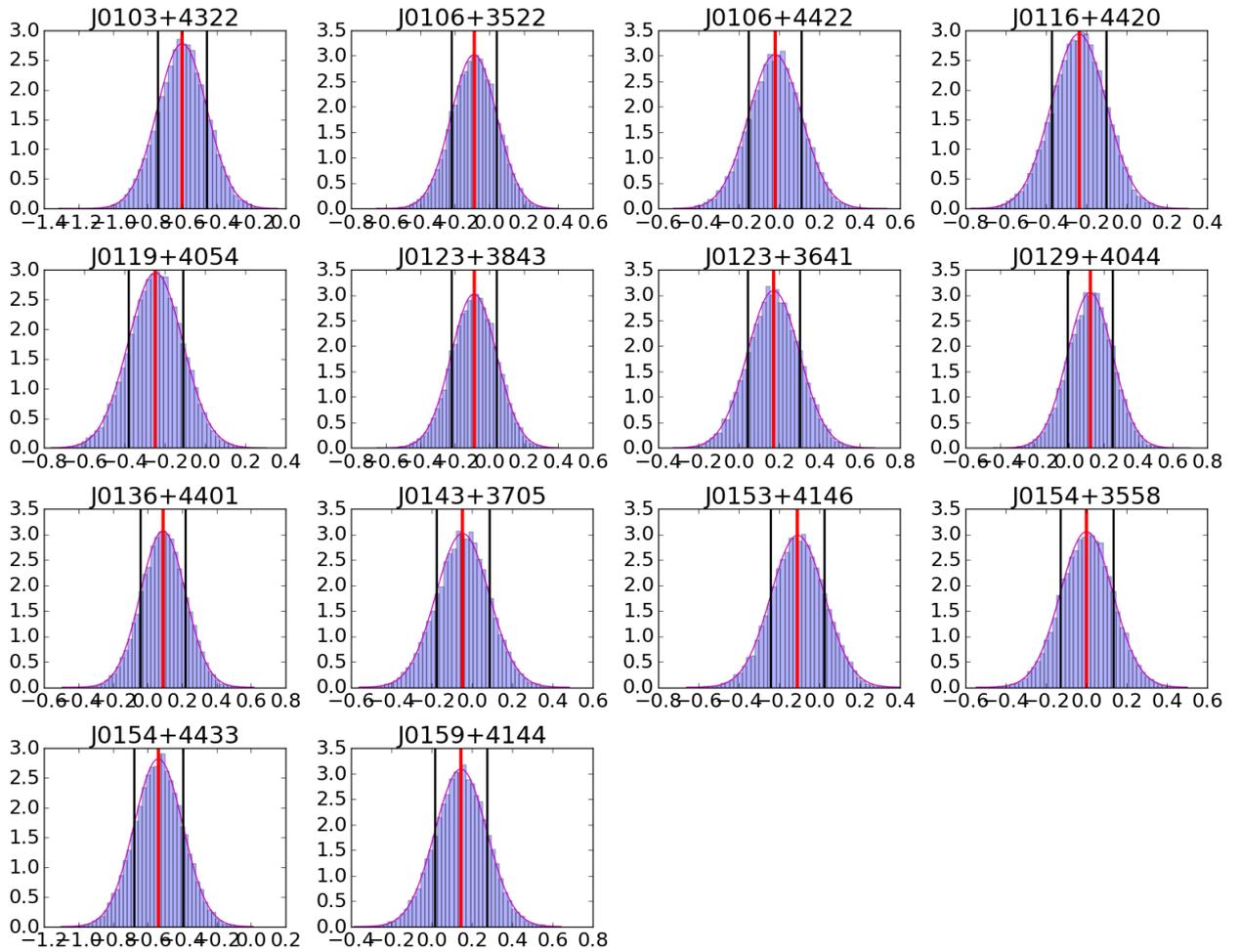


Figure 21: The log-redshift distributions around the sample K-band measurements. The red line indicates the mean, the black lines the symmetric dispersion. The red curve is the fitted continuous Gaussian distribution through the histogram data. The x-axes give the  $\log_{10}(z)$ -value, the y-axes the probability in arbitrary units.

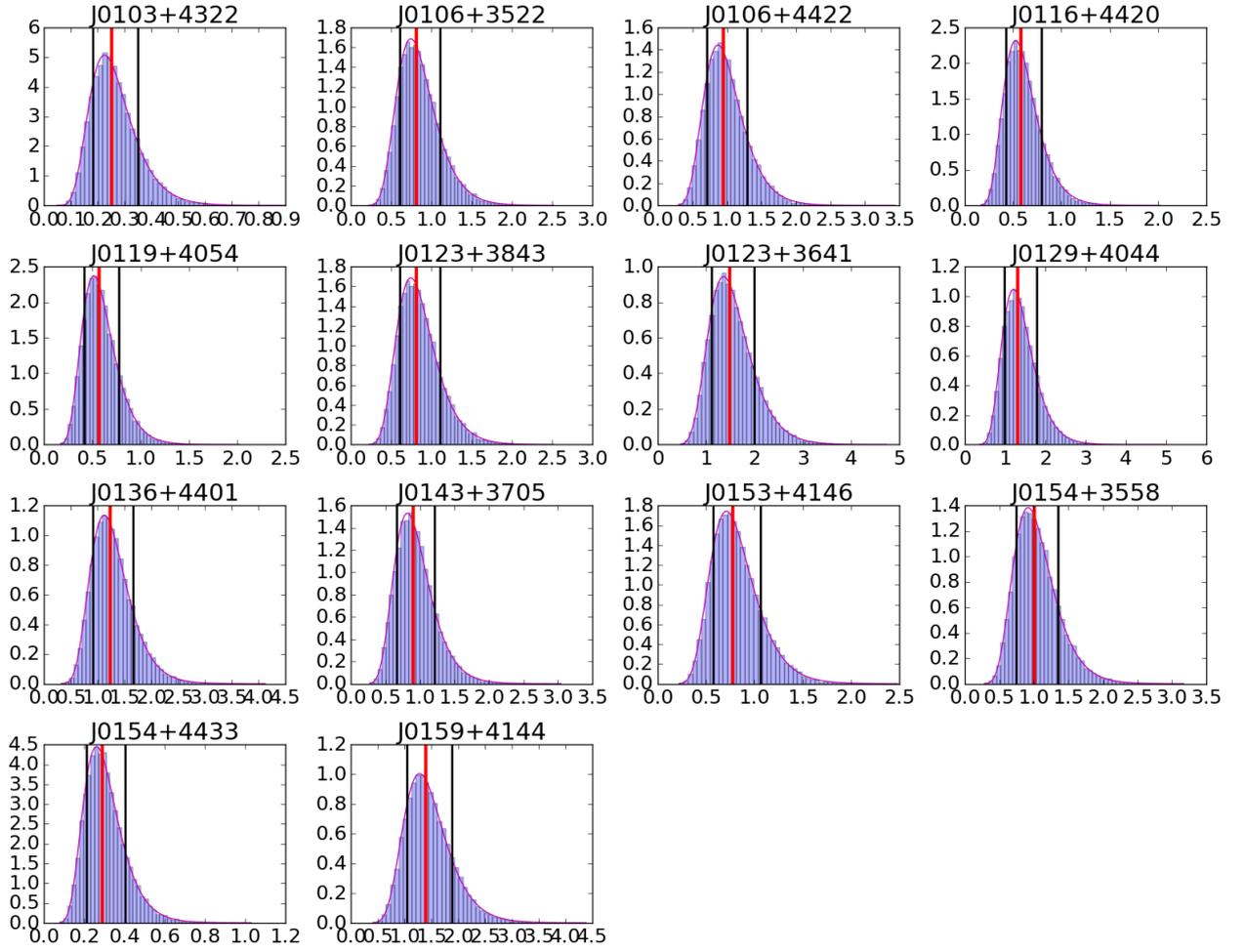
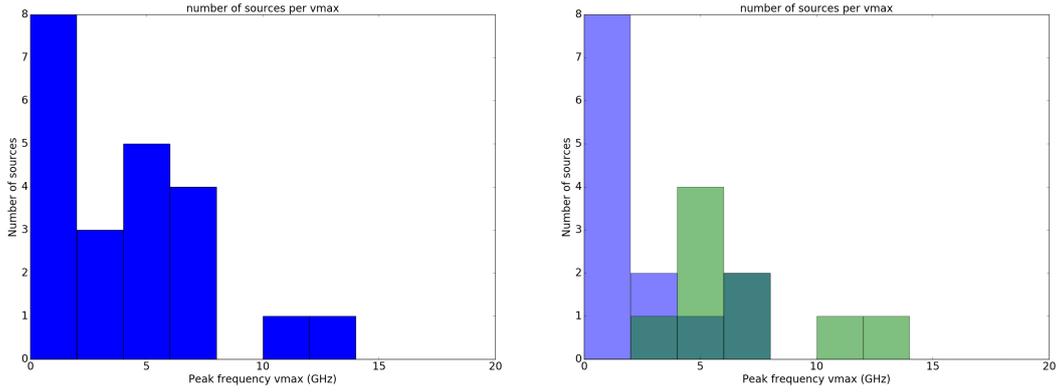
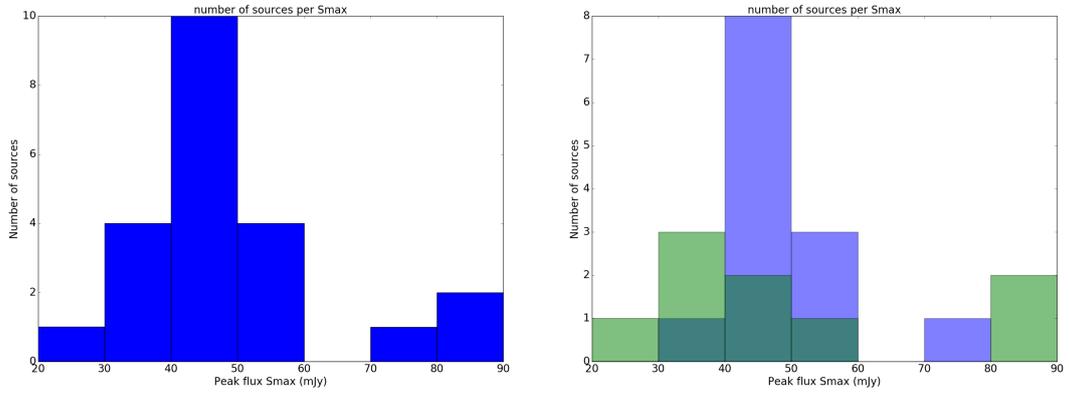


Figure 22: The redshift distributions around the sample K-band measurements. The red line indicates the median, the black lines the lower and upper boundaries of the 68% confidence interval. The red curve is the fitted continuous log-normal distribution through the histogram data. The x-axes give the  $z$ -value, the y-axes the probability in arbitrary units.



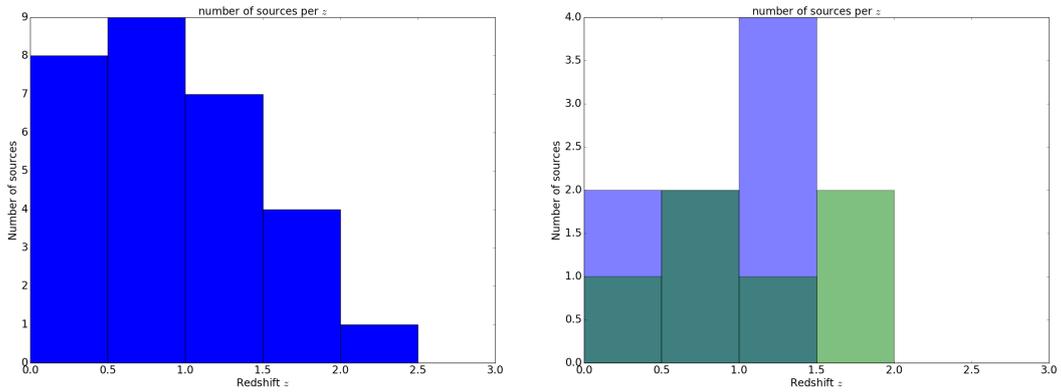
(a) SSA GPS and SSA+power law GPS populations combined. (b) Separate GPS populations (blue: SSA, green: SSA+power law).

Figure 23: Number counts for peak frequency  $\nu_{max}$ . Bin width is 2 GHz.



(a) SSA GPS and SSA+power law GPS populations combined. (b) Separate GPS populations (blue: SSA, green: SSA+power law).

Figure 24: Number counts for peak flux density  $S_{max}$ . Bin width is 10 mJy



(a) SSA GPS and SSA+power law GPS populations combined. (b) Separate GPS populations (blue: SSA, green: SSA+power law).

Figure 25: Number counts for redshift  $z$ . Bin width is 0.5 redshift.

## 4.4 Results

The redshift results for the sample that underwent the K-z treatment are provided in the upper section of Table 7. Combined with the rest of the sample, with spectroscopic redshifts taken from [Marlow et al., 2000] and NASA Extragalactic Database (NED), we obtain the full table. These values can be translated to a histogram of number count of sources per redshift bin (Figure 25). This has also been done for peak frequency (Figure 23) and peak flux (Figure 24) of the GPS and GPS+power law sources (see Appendix B.5). While the separate populations of SSA GPS and SSA+power law GPS in the peak-related Figures add up to their combined counterparts, in Figure 25 this is not the case. This is because the combined plot also includes power law spectra whose redshift is known.

When studying the redshift number counts it becomes apparent that there is no clear sign of bimodal behaviour regarding the two separate populations. It was suggested by Dr. McKean that a bimodal trend might arise between a population at small redshift (radio galaxies) and one at large redshift (radio quasars), possibly coinciding with the separation between SSA GPS and SSA+power law GPS sources. Figure 25b however clearly shows that this is not the case, and from Figure 25a we can continue to conclude that there is no secondary peak at higher redshift.

A slightly more differentiated situation can be spotted in Figures 23 and 24. In frequency space, it can be observed that the peak frequency of pure SSA GPS sources is in general lower than those of their counterparts in the SSA+power law GPS sample, with an exceptional spike between 0 and 2 GHz. The SSA+power law GPS population shows a preference towards the multi-GHz regime, with sporadic counts exceeding 10 GHz prompting the suggestion that these would be rather classified as HFPs. In terms of peak flux, another trend becomes apparent with the SSA+power law GPS possessing an on average lower peak flux density than their pure SSA GPS counterparts.

Combining all redshifts obtained by using the K-z relation with spectroscopic redshifts from NED and Marlow et al. (2000) and averaging their squares, then taking the root of this quantity, we obtain a root mean square (RMS) redshift - the RMS is a useful statistic which gives an impression of the spread of a quantity in its space. Marlow et al. obtained an RMS of 1.27 with spread of 0.95 for a completeness level of 64% [Marlow et al., 2000] for the CLASS sample. It is interesting to note, however, that these values were seemingly obtained by setting those redshift values that Marlow et al. failed to determine spectroscopically to 0, a practice that this thesis does not follow. Hence the RMS was determined from the raw Marlow redshift data, which for the sources in Marlow's sample with known redshifts gave RMS  $\langle z \rangle = 1.77$  with an RMS spread of 1.01, rather than the value mentioned in the paper's abstract. Combining the Marlow et al. CLASS sample with the sources used in this thesis, an RMS of  $\langle z \rangle = 1.45$  with an RMS spread of 0.87 for a completeness level of 66% was obtained. Sources within this thesis' sample with unknown redshifts were simply omitted from this computation, rather than taken at  $z = 0$ . This shows that the redshifts determined by this paper using the K-z relation significantly drift the RMS of the sample towards lower redshifts, indicating that McKean's observed sources are on average closer than Marlow's CLASS sample.

## 5 Discussion

This chapter discusses the synthesis of the results so far obtained. Cosmological statistics of the sample will be provided by combining the results of the spectral fitting and the K-z relation, an analysis and discussion will be provided regarding these final results and suggestions will be offered for possible next steps.

### 5.1 Synthesis of results

Comparing the results for the spectral fitting and the determination of redshift, be it by spectroscopy or by using the K-z relation, a sub-sample was constructed of those sources with a spectral turnover. In case a redshift was not known, the mean value of the other redshifts ( $\sim 1.01$ ) was used as an informed estimate.

Source	redshift $z$	$S_{max}$ (mJy)	$\nu_{obs}$ (GHz)	$\theta$ (mas)	$D$ (Mpc)	$d$ (pc)	$\nu_e$ (GHz)
J0104+3840 (!)	$1.00^{+0.35}_{-0.26}$	$45.3 \pm 18.6$	$0.79 \pm 0.02$	$0.28^{+2.60}_{-0.28}$	6879.0	2.3	1.58
J0106+3522 (!)	$0.82^{+0.29}_{-0.21}$	$41.3 \pm 2.1$	$1.13 \pm 0.02$	$0.36^{+0.39}_{-0.33}$	5347.5	2.8	2.05
J0106+4422 (!)	$0.95^{+0.34}_{-0.25}$	$73.7 \pm 6.1$	$1.20 \pm 0.02$	$0.4^{+0.59}_{-0.40}$	6496.0	3.3	2.34
J0123+3641 (!)	$1.48^{+0.51}_{-0.38}$	$55.9 \pm 1.7$	$1.99 \pm 0.02$	$0.22^{+0.25}_{-0.19}$	11201.0	1.9	4.94
J0128+4439	$0.228 \pm 0.001$	$40.3 \pm 0.4$	$2.88 \pm 0.02$	$2.02^{+0.02}_{-0.02}$	1174.9	7.6	3.54
J0154+3558	$1.01^{+0.35}_{-0.26}$	$35.9 \pm 4.3$	$3.08 \pm 0.02$	$0.3^{+0.15}_{-0.14}$	6947.4	2.5	6.19
J0154+4433	$0.29^{+0.11}_{-0.08}$	$48.8 \pm 1.6$	$5.01 \pm 0.02$	$1.83^{+0.03}_{-0.03}$	1553.0	8.3	6.47
J0159+4144 (!)	$1.40^{+0.48}_{-0.36}$	$54.1 \pm 8.1$	$1.86 \pm 0.02$	$0.23^{+0.44}_{-0.23}$	10426.4	2.0	4.45
J1702+5511 (!)	$1.00^{+0.35}_{-0.26}$	$40.7 \pm 3.6$	$0.90 \pm 0.02$	$0.27^{+0.69}_{-0.27}$	6879.0	2.2	1.80
J1702+5812 (!)	$1.00^{+0.35}_{-0.26}$	$42.0 \pm 0.6$	$1.95 \pm 0.02$	$0.30^{+0.18}_{-0.14}$	6879.0	2.5	3.89
J1726+5826 (!)	$1.00^{+0.35}_{-0.26}$	$49.1 \pm 0.6$	$1.04 \pm 0.02$	$0.30^{+0.48}_{-0.30}$	6879.0	2.5	2.08
J1728+5532	$1.404 \pm 0.002$	$42.0 \pm 1.4$	$7.17 \pm 0.02$	$0.26^{+0.01}_{-0.01}$	10486.6	2.3	17.25
J1758+5951	$1.00^{+0.35}_{-0.26}$	$59.8 \pm 5.2$	$7.26 \pm 0.02$	$0.47^{+0.05}_{-0.05}$	6879.0	3.9	14.53
J0103+4322	$0.25^{+0.10}_{-0.07}$	$49.2 \pm 4.4$	$12.07 \pm 0.01$	$2.69^{+0.02}_{-0.02}$	1306.2	10.9	15.09
J0123+3843	$0.82^{+0.29}_{-0.21}$	$36.2 \pm 0.5$	$5.80 \pm 0.02$	$0.45^{+0.04}_{-0.03}$	5347.5	3.5	10.53
J0123+3806	$1.656 \pm 0.004$	$89.6 \pm 1.0$	$6.20 \pm 0.02$	$0.3^{+0.01}_{-0.01}$	12866.8	2.6	16.46
J0131+4428	$1.123 \pm 0.001$	$38.2 \pm 0.9$	$11.77 \pm 0.02$	$0.36^{+0.01}_{-0.01}$	7945.6	3.1	24.98
J0143+3705	$0.90^{+0.32}_{-0.24}$	$53.7 \pm 1.1$	$4.81 \pm 0.02$	$0.47^{+0.06}_{-0.05}$	5995.4	3.8	9.11
J0151+4417	$1.976 \pm 0.003$	$85.6 \pm 0.9$	$4.73 \pm 0.02$	$0.22^{+0.01}_{-0.01}$	15999.0	1.9	14.07
J0159+4059	$1.00^{+0.35}_{-0.26}$	$40.1 \pm 1.5$	$3.29 \pm 0.02$	$0.33^{+0.09}_{-0.07}$	6879.0	2.7	6.59
J1720+5541	$1.00^{+0.35}_{-0.26}$	$26.0 \pm 1.4$	$5.18 \pm 0.02$	$0.29^{+0.03}_{-0.02}$	6879.0	2.4	10.37
J1746+5659	$1.00^{+0.35}_{-0.26}$	$35.0 \pm 0.2$	$7.49 \pm 0.02$	$0.36^{+0.01}_{-0.01}$	6879.0	3.0	14.98

Table 8: Cosmological statistics for those 22 sources in the sample with peaks. All objects with redshift  $1.00^{+0.35}_{-0.26}$  had their redshifts estimated by taking the approximate average of all the known redshifts in the sample, namely (very close to) unity. All symbols take their usual meanings, while  $\theta$  denotes the angular extension of the source (angular diameter in the sky),  $D$  the luminosity distance to the source,  $d$  the corresponding linear extension and  $\nu_e$  the emitted frequency of the spectral peak. Above the horizontal line are SSA GPS; below are SSA GPS+power law spectra. Those SSA GPS sources marked with (!) deviate from the expected trend in Figure 26.

#### 5.1.1 Angular size

Given the peak flux, the redshift, the angular size  $\theta$  and the magnetic field strength  $B$ , the peak frequency of a source can be computed via,

$$\nu_{max} = f(k) \cdot B^{1/5} S_{max}^{2/5} \theta^{-4/5} (1+z)^{1/5}. \quad (38)$$

[Kellermann and Pauliny-Toth, 1981] This can be rewritten for the angular size as,

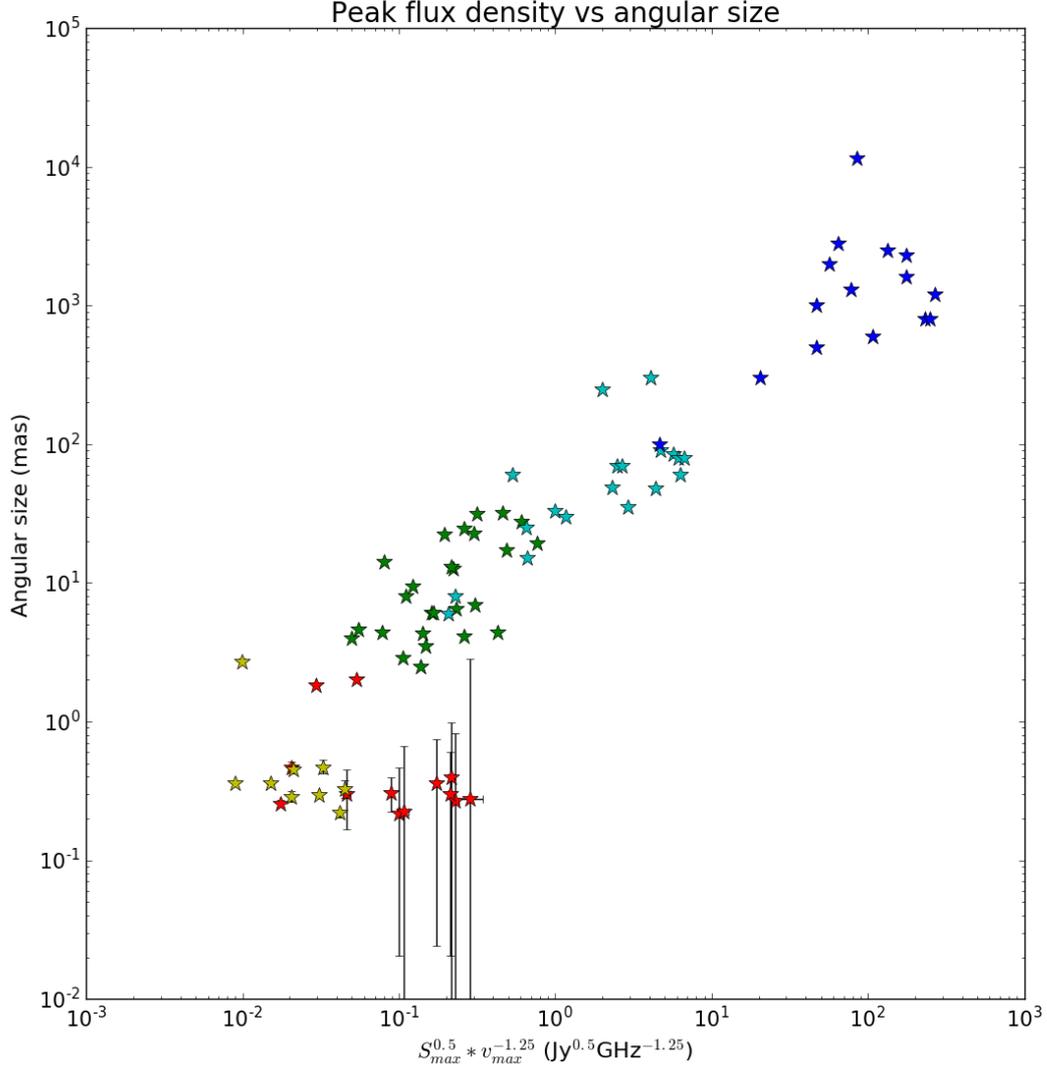


Figure 26: Angular size of source versus  $S_{max}^{0.5} \nu_{max}^{-1.25}$ . This plot is made within the observers frame. Cyan, blue and green stars are taken from [Snellen et al., 2000a], and designate Snellen’s CSS, bright GPS and faint GPS samples respectively. The red stars represent the sources of McKean’s very faint SSA GPS sample, while the yellow stars stand for the SSA GPS+power law sources found by this thesis.

$$\theta = f(k)^{5/4} \cdot \frac{B^{1/4} S_{max}^{1/2} (1+z)^{1/4}}{\nu_{max}^{5/4}}. \quad (39)$$

Wherein  $\nu_{max}$  is the GPS peak frequency,  $f(k)$  a function that depends weakly on the thick spectral index, but which is assumed as approximately 8 (the value it takes for  $k \sim 2$ ),  $B$  the magnetic field strength,  $S_{max}$  the GPS peak flux density,  $\theta$  the angular size and  $z$  the redshift. Furthermore,  $B$  is taken as  $50 \mu\text{G}$ , based on Mingaliev et al. (2013) who used  $100 \mu\text{G}$  for their sample of CRSs: however, as our sample contains fainter sources than those in Mingaliev’s sample and thus sources with less synchrotron emission which indicates the presence of a weaker magnetic field, we take a smaller magnetic field strength -  $50 \mu\text{G}$  was chosen in discussion with Dr. McKean as a sensible value. This led to the computation

$S_{med}$ (mJy)	$S_{bin}$ (mJy)	n	$\frac{dN}{dS} S_{med}^{5/2}$ ( $\text{Sr}^{-1}\text{Jy}^{3/2}$ )	$\frac{dN}{dS}$ ( $\text{Sr}^{-1}\text{Jy}^{-1}$ )
30	10.0	5	$0.83 \pm 0.37$	$(16.74 \pm 7.49) \cdot 10^{-5}$
60	20.0	15	$7.00 \pm 1.81$	$(25.11 \pm 6.48) \cdot 10^{-5}$
60	20.0	7 (8)	$3.27 \pm 1.24$	$(11.72 \pm 4.43) \cdot 10^{-5}$
120	40.0	2	$2.64 \pm 1.87$	$(1.67 \pm 1.18) \cdot 10^{-5}$

Table 9: The values of the bins for McKean’s sample.  $S_{med}$  is the average bin flux density,  $S_{bin}$  the bin’s radius, n the plain number of sources per bin (in brackets on the third row is provided the number of sources which are offset from the trend in Figure 26, which are all in the 60 mJy bin),  $\frac{dN}{dS}$  the number of sources per sky area per bin and  $\frac{dN}{dS} S_{med}^{5/2}$  that number density corrected by the average bin flux density. Higher bins did not contain any of the McKean sources and were thus omitted in this table. The second row is duplicated, once taking into account all sources, and secondly without those sources offset from the trend in Figure 26

$S_{med}$ (mJy)	$S_{bin}$ (mJy)	n	$\frac{dN}{dS} S_{med}^{5/2}$ ( $\text{Sr}^{-1}\text{Jy}^{3/2}$ )	$\frac{dN}{dS}$ ( $\text{Sr}^{-1}\text{Jy}^{-1}$ )
30	10.0	3	$0.08 \pm 0.05$	$(1.62 \pm 0.93) \cdot 10^{-5}$
60	20.0	7	$0.53 \pm 0.20$	$(1.89 \pm 0.71) \cdot 10^{-5}$
120	40.0	8	$1.70 \pm 0.60$	$(1.08 \pm 0.38) \cdot 10^{-5}$
240	80.0	12	$7.23 \pm 2.08$	$(8.09 \pm 2.34) \cdot 10^{-6}$
480	160.0	1	$1.70 \pm 1.70$	$(3.37 \pm 3.37) \cdot 10^{-7}$

Table 10: The values of the bins for Snellen’s sample. The columns represent the same quantities as in Table 9.

of angular size for our sub-sample, which turned out to be in the milli-arcseconds regime as one would expect for GPS sources. Further statistics (luminosity distance  $D$ , linear extension  $d$  and peak frequency corrected for redshift  $\nu_{emit}$ ) were computed using respectively the source code of a web-based Cosmology Calculator [Wright, 2006] and equation 31, which were all combined in Table 8. As was done in Snellen et al. (2000b), angular size was then plotted against a peak convolution given by  $S_{max}^{0.5} \nu_{max}^{-1.25}$ , results of which are shown in Figure 26 (see Appendix B.6).

The majority of the McKean sources line up well with Snellen’s various samples, although a subset of the sources analysed in this thesis are offset from the trend. These coincidentally also possess large uncertainty bars in their angular size, while their uncertainty bars in the x-direction do not allow them to fall within the regime of the other, better fitting sources. There is a strong correspondence between offset sources and the spectral type: there are no SSA GPS+power law sources in the offset subset. There also seems to be a correspondence between the large uncertainties and poor spectral fits, although this is not exclusively the case. This can either indicate that indeed in the fitting procedure some parameter or function was poorly defined, or that these sources represent a new population of GPS that are not well understood. The 8 sources which both possess large uncertainties and are offset from the trend are J0104+3840, J0106+3522, J0106+4422, J0123+3641, J0159+4144, J1702+5511, J1702+5812 (although this source is a limit case, having uncertainties that are small compared to the other 7) and J1726+5826. These are exclusively SSA GPS sources with peak frequencies around 1 GHz, while the SSA GPS+power law sources line up consistently better with the trend provided by Snellen’s sources. Further inspection shows that those SSA GPS sources that do line up with the trend all have peak frequencies in excess of  $\sim 2$  GHz. This leads us to conclude that the correspondence between peak frequency and angular size might be different than originally suspected in the low flux-density and small angular size regime. Another option is that the sources with low-frequency peaks suffer major variability: as the high- and low-frequency observations were not carried out simultaneously, it is well possible that if they were repeated in one sitting, the spectra would look markedly different, possibly shifting the peak to higher regimes.

$S_{med}$ (mJy)	$S_{bin}$ (mJy)	n	$\frac{dN}{dS} S_{med}^{5/2}$ ( $\text{Sr}^{-1}\text{Jy}^{3/2}$ )	$\frac{dN}{dS}$ ( $\text{Sr}^{-1}\text{Jy}^{-1}$ )
30	10.0	8	$0.91 \pm 0.32$	$(18.36 \pm 6.49) \cdot 10^{-5}$
60	20.0	22	$7.53 \pm 1.61$	$(27.00 \pm 5.76) \cdot 10^{-5}$
120	40.0	10	$4.34 \pm 1.37$	$(2.75 \pm 0.87) \cdot 10^{-5}$
240	80.0	12	$7.22 \pm 2.08$	$(8.09 \pm 2.33) \cdot 10^{-6}$
480	160.0	1	$1.70 \pm 1.70$	$(3.37 \pm 3.37) \cdot 10^{-7}$

Table 11: The values of the bins for the combined sample. The columns represent the same quantities as in Table 9.

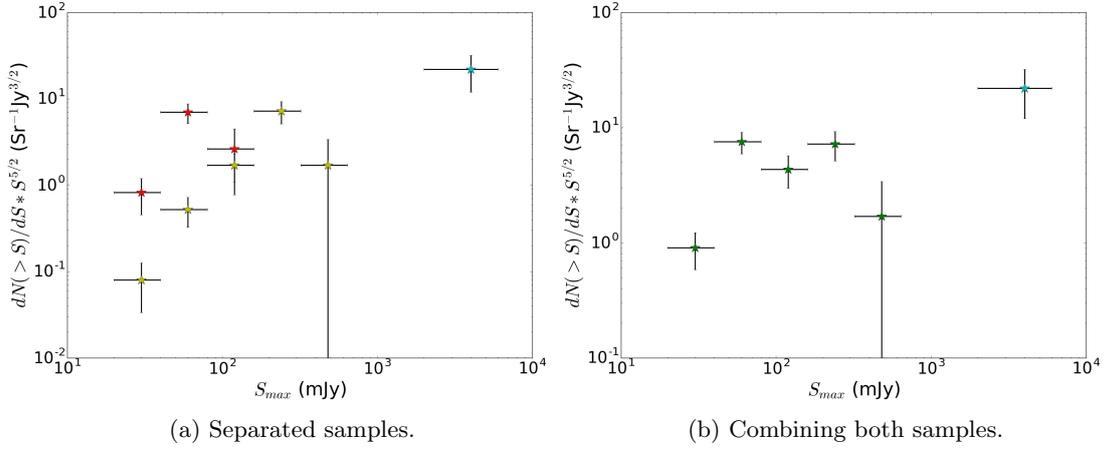


Figure 27: Differential number counts/peak convolution versus the peak flux. Red stars are the McKean sources, yellow stars are Snellen's, the cyan star was derived by Snellen from [de Vries et al., 1997]. The green stars represent the combined sample (bin width indicated by horizontal bar through stars).

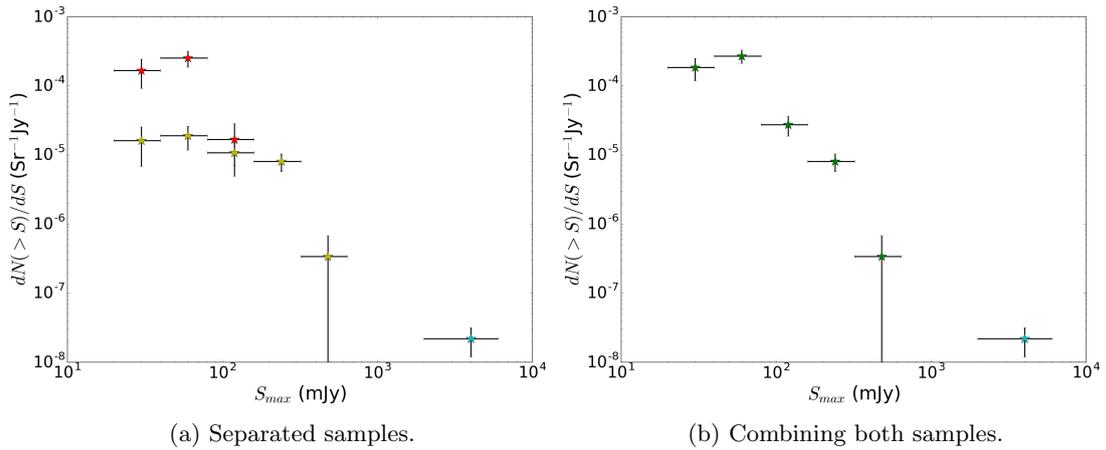
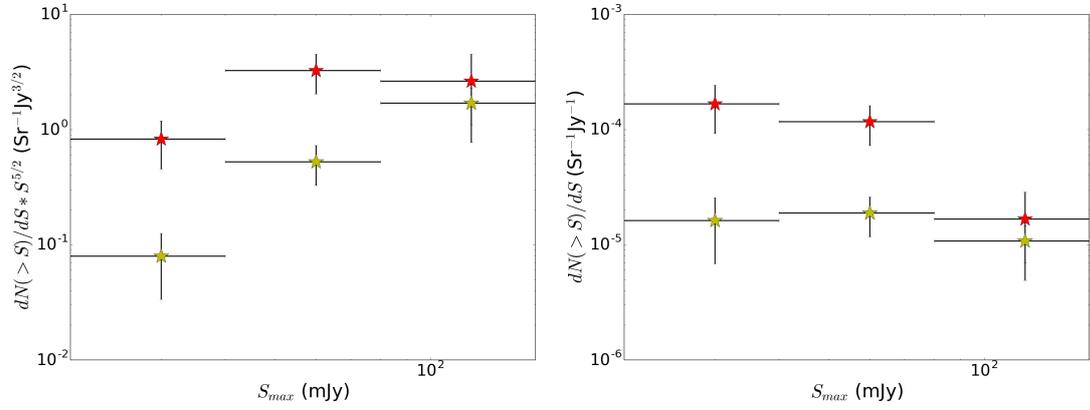
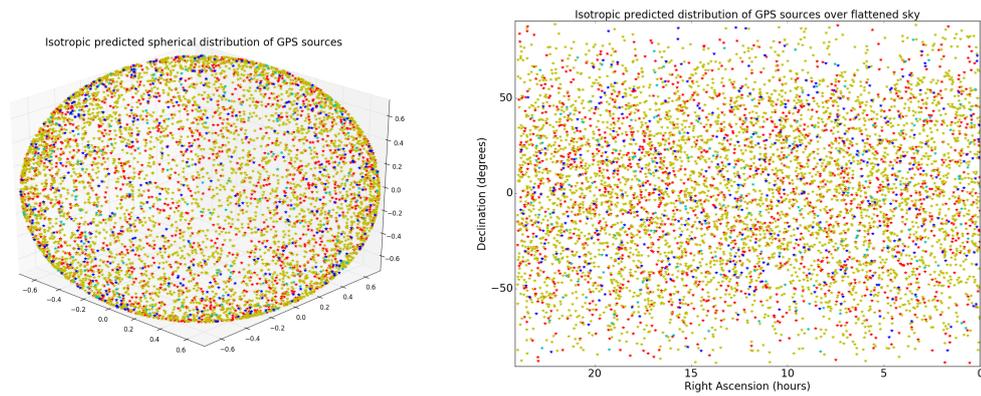


Figure 28: Plain differential number counts versus the peak flux. Red stars are the McKean sources, yellow stars are Snellen's, the cyan star was derived by Snellen from [de Vries et al., 1997]. The green stars represent the combined sample (bin width indicated by horizontal bar through stars).



(a) Discrepancy in number count/peak convolution versus peak flux. (b) Discrepancy plain differential number count versus peak flux.

Figure 29: Closeup of the discrepancy between the McKean and Snellen samples, after removing those sources offset from the trend in Figure 26. Red stars are the McKean sources, yellow stars are Snellen's (bin width indicated by horizontal bar through stars).



(a) Populations projected onto an orb. (b) Populations projected onto a plain.

Figure 30: Predicted population densities of GPS. The red population is 20-40 mJy, yellow 40-60 mJy, cyan 60-80 mJy and blue >80 mJy. Note that the locations are completely arbitrary, and these plots serve only to give an impression of the observed density of these sources.

### 5.1.2 Number density per peak flux density

Snellen, in his 1998 paper, creates a plot of differential number count of GPS sources, i.e. the number of sources within some bin of peak flux density multiplied by the average flux of the bin to the power 5/2 [Snellen et al., 1998b]. Using the data acquired by this thesis, an expanded version of this plot was created. Firstly, to obtain compatible density values, sky areas were computed for both the Snellen et al. (1998) and the McKean sample, using equation 24. This yielded a sky area for Snellen of 0.293 steradians, while the McKean field extended 0.047 steradians. Snellen’s sources were subjected to the same criteria used to select the McKean sample (see chapter 3, section 1), namely a spectral index of no steeper than -0.5 between 1.4 and 4.85 GHz and a flux density of no smaller than 16.7 mJy at 8.46 GHz. These sources were then combined with McKean’s and divided into bins (which were maintained of constant size in  $\log_{10}$ -space) dependent on peak flux density to create Figures 27 and 28 (see Appendix B.7). The bin values, both in terms of number of sources, and the plotted values before and after multiplication by  $S_{max}^{5/2}$ , are presented in Tables 9, 10 and 11.

These Figures show a clear deviation from the trend expected from the Snellen and De Vries data alone. Even when one does not correct the Snellen data via the selection criteria set by McKean, resulting in more sources in the Snellen bins, his data still does not connect directly to the McKean sample and a discrepancy between Snellen’s lowest and McKean’s highest entry remains. This contradicts expectations, as what one would intuitively expect to see is the populations to i) link up directly to make a continuous curve and ii) the population at low flux density to shrink. This could hint at an unexplored, large population of low-flux density GPS, but it can also be due to, as previously mentioned, sources which are highly variable and which, due to non-simultaneous measurements of their high- and low-frequency regimes, show falsely peaked spectra and as such should therefore not even be included in these number density graphics. Further investigation (Figure 29) however shows that even after the removal of those possibly non-peaked variable sources masquerading as GPS, there is a considerable discrepancy of almost an order of magnitude at most between the McKean and Snellen samples, necessitating another explanation for the remaining difference.

One can use the number density algorithm developed for these plots to make a prediction of the amount of sources within some flux bin in the night sky. This is displayed in Figure 30. The positions of these sources are arbitrary, but the density impression that it gives is found by McKean’s data, further giving us a predicted amount of sources satisfying McKean’s criteria of approximately  $\sim 13000$ . The 40-60 mJy population clearly dominates (see Appendix B.4).

## 5.2 Conclusions regarding the evolution of radio sources

Combining all results so far obtained, the following can be stated about this sample of very faint GPS sources: that they are small, both in angular ( $\sim 1$  mas) and linear (several pc) terms, and that their emitted peaks are generally within the multi-GHz regime. We can also state that SSA GPS+power law sources in general have higher peak frequencies, which is sensible taking into account that the power law dominates the lower frequency regions of the spectrum, but there does not seem to be either a linear extension nor a redshift division between the two populations. In addition, we can conclude that during their evolution from very faint GPS (McKean’s sample) to bright CSS (Snellen’s sample), the relation between angular extension and  $S_{max}^{0.5} \cdot \nu_{max}^{-1.25}$  remains approximately constant (see Figure 39), although an interesting subsection of pure SSA GPS sources seem offset from this trend. This implies that pure GPS sources that have just been energized increase in angular size quite explosively in the first stage of their evolution, by approximately an order of magnitude, before suffering changes in their peak flux density or peak frequency and joining the trend set out by Snellen’s sample. It is also possible that during this early phase,  $S_{max}^{0.5}$  and  $\nu_{max}^{1.25}$  change inversely proportionally, i.e. the flux goes up as the frequency, as predicted by CRS evolution models, goes down, leading to an approximately constant quotient over this period. This is somewhat unexpected as SSA spectra ordinarily see both their flux *and* their frequency go down with time. However, the mechanism behind this change, if at all physical, is unknown at the time. It is also possible that the relation between angular size and peak flux and frequency in fact changes as the source ages, or that those offset sources are in fact non-peaked but highly variable power law sources masquerading as GPS. However, analysis done by making number counts per flux bin shows that even after accounting for this possible issue, there is still a large discrepancy between McKean’s and Snellen’s samples, raising questions concerning whether there might be a larger than expected population of low

flux density GPS in the night sky.

This roughly constant behaviour raises suspicions that these sources are indeed of the same nature, and that, as first hypothesised by O’Dea, Baum and Stanghellini, GPS do in fact over time evolve into CSS, although further research is required before a definitive conclusion can be reached, particularly into the transition class between CSS and FRI/II radio galaxies. Despite that, the findings so far presented provide a tantalizing first glimpse into the very early evolution of cosmic radio sources. In addition, the discovery of a dense population of low flux-density sources presents an unexpected upturn in the number count in that regime and will need to be investigated thoroughly in future research, as these objects may hold the clue to what happens in the earliest days of the life of an AGN. That there are so much more of these faint objects than brighter ones might imply that not all GPS sources go on to become bright radio galaxies, instead fading far earlier.

### 5.3 Next steps

Research is never complete, and this open scientific frontier forms no exception. Advances in this field can be enabled by several actions. Above all else, observations of both the very faint GPS - possibly even the HFP - and the CSS-to-FRI/II transition populations would greatly enhance our understanding of the evolution of cosmic radio sources. The actual triggering of CRSs still remains largely an enigma, as does the physical transition stage from CSS to a fully-fledged radio galaxy. Observations of these populations would enable us to study these critical evolutionary stages in unprecedented detail, and although the McKean sample forms a revolutionary, pioneering first step into the very faint CRS regime, it is hoped that it is not by far the last.

Additional improvements to our conclusions could be made with additional, more detailed observations of sources within the Snellen flux density observation window ( $\sim 40$  to  $\sim 900$  mJy). His observations stem from 1998, and although they are highly important and his paper ([Snellen et al., 1998b]) has been more than critical to the completion of this project, until a thorough re-examination of the moderate GPS population is performed, the option cannot be ruled out that the reason behind the discrepancy in Figures 27 and 28 is simply due to an incompleteness in the higher flux density samples.

The quality of the spectral fits could also be greatly improved by additional measurements in the observed band (0.325 - 22.46 GHz), especially on the low end of this range. The majority of fitting issues stemmed from the fact that either  $k$  (in SSA GPS sources) or  $m$  (in SSA GPS+power law sources) was poorly constrained, often by just two parameters. The upgraded VLA should be able to perform these observations, which would undoubtedly monumentally increase the quality of the fits, constraining the parameters more strongly than the six data points given ever could. It would also improve our fits if the measurements at 325 MHz and 1.4 GHz were redone simultaneously with the VLA measurements, as the current observations at the low and high ends of the spectrum lie years apart, something which presents an issue given the variable nature of some GPS sources. In the end, most uncertainties in the final results can be traced back to uncertainties within the spectral fits, so if they were to be more properly constrained and informed, so would all following physical results. Finally, spectroscopic analysis of those sources in the sample that lacked any redshift would greatly improve the completeness level of the sample, and would shed more light on the nature and shape of the redshift number counts, proving definitively the existence or non-existence of a bimodal behaviour in terms of redshift between the SSA GPS and SSA GPS+power law populations. Essentially, more data would be greatly appreciated.

### 5.4 Summary and concluding remarks

In short: a sample of 45 compact radio sources have been fitted with SSA spectra and/or power laws. The ratio between peaked and non-peaked spectra proved approximately 1:1, with most peaks in the multi-GHz regime. Redshifts of a subsample were determined using the K-z relation, while others were found to have spectroscopic redshifts, resulting in a completeness level of 66.6%. This data was combined to obtain the following conclusions:

1. there is no sign of bimodal behaviour in terms of redshift number counts between SSA GPS and SSA GPS+power law spectrum sources;
2. there are weak signs of bimodal behaviour in terms of peak flux and peak frequency number counts between SSA GPS and SSA GPS+power law spectrum sources;

3. the angular sizes of sources within the sample were of the scale  $\sim 0.5$  mas;
4. the linear sizes of sources within the sample were of the scale  $\sim 3$  pc;
5. the relation between angular size and  $S_{max}^{0.5} \cdot \nu_{max}^{1.25}$  remains linear at low flux densities for SSA GPS+power law sources and falls off sharply at low flux densities for SSA GPS sources with respect to sources at larger flux densities, as obtained by Snellen, for which the origin or mechanism is unknown;
6. there is an unexpectedly dense population of low flux density sources present, compared to higher flux density, which may indicate that not all GPS sources become CSS or FRI/II radio galaxies.

In the end, suggestions are given on how to improve future research. These include measurements at both the extreme high and low frequency ends of the presently known CRS population, a redo of the sources within Snellen's observed flux density window, and additional observations at different frequencies of the sources in this sample to improve spectral fits.

# Appendices

## A GPS sample observations

These tables have been taken from [McKean, Browne & Jackson, in prep.]. They list the observations of the GPS sample used in this thesis. Note that in the second table, the source J0112+3759 is missing.

**Table A1.** The sample positional information with 1.4, 4.85 and 8.46 GHz radio flux-densities measured from NVSS, GB6 and CLASS, respectively.

Source Name	Right Ascension (° ' ")	Declination (° ' ")	$S_{0.325}$ (mJy)	$S_{1.4}$ (mJy)	$S_{4.85}$ (mJy)	$S_{8.46}$ (mJy)	$\alpha_{1.4}^{4.85}$
J0103+4322	01 03 28.8025	+43 22 59.526	45.0 ± 3.7	44.6 ± 1.6	41.0 ± 5.0	37.0 ± 1.9	-0.07 ± 0.10
J0104+3840	01 04 22.2100	+38 40 35.400	< 18.0	40.5 ± 2.7	27.0 ± 4.0		-0.33 ± 0.13
J0105+3533	01 05 57.2500	+35 33 29.400	66.0 ± 3.4	43.0 ± 1.3	28.0 ± 4.0		-0.35 ± 0.12
J0106+3522	01 06 14.4844	+35 22 26.786	< 18.0	40.5 ± 1.3	29.0 ± 5.0	18.5 ± 0.9	-0.27 ± 0.14
J0106+4422	01 06 21.3985	+44 22 27.377	33.0 ± 4.1	72.8 ± 2.2	47.0 ± 5.0	32.5 ± 1.6	-0.35 ± 0.09
J0112+3759	01 12 00.2598	+37 59 31.559	143.0 ± 3.7	61.9 ± 1.9	37.0 ± 5.0	21.8 ± 1.1	-0.41 ± 0.11
J0116+4420	01 16 56.8405	+44 20 59.871	169.0 ± 4.2	69.6 ± 2.5	41.0 ± 5.0	28.8 ± 1.5	-0.43 ± 0.10
J0119+4054	01 19 47.9105	+40 54 18.586	86.0 ± 4.0	75.2 ± 2.8	45.0 ± 5.0	48.4 ± 2.4	-0.41 ± 0.09
J0123+3843	01 23 23.8308	+38 43 57.900	< 18.0	18.5 ± 0.7	27.0 ± 4.0	36.6 ± 1.8	+0.30 ± 0.12
J0123+3806	01 23 28.8725	+38 06 36.758	46.0 ± 3.8	36.8 ± 1.5	40.0 ± 5.0	63.4 ± 3.2	+0.07 ± 0.11
J0123+3641	01 23 58.5880	+36 41 45.336	35.0 ± 3.6	54.0 ± 2.0	31.0 ± 4.0	28.9 ± 1.5	-0.45 ± 0.11
J0128+4003	01 28 13.6382	+40 03 29.866	116.0 ± 3.8	56.3 ± 1.7	47.0 ± 5.0	28.9 ± 1.5	-0.15 ± 0.09
J0128+4439	01 28 41.3396	+44 39 17.982	< 18.0	36.9 ± 1.2	35.0 ± 5.0	60.8 ± 3.0	-0.04 ± 0.12
J0129+4044	01 29 59.3005	+40 44 02.981	175.0 ± 3.6	67.9 ± 2.1	38.0 ± 5.0	24.8 ± 1.3	-0.47 ± 0.11
J0131+4428	01 31 03.3632	+44 28 01.883	65.0 ± 3.6	33.8 ± 1.4	27.0 ± 4.0	24.4 ± 1.5	-0.18 ± 0.12
J0132+4345	01 32 06.0642	+43 45 34.428	98.0 ± 3.7	55.0 ± 1.7	41.0 ± 5.0	30.1 ± 1.5	-0.24 ± 0.10
J0134+4018	01 34 40.8057	+40 18 18.619	104.0 ± 3.6	50.7 ± 1.6	29.0 ± 4.0	19.6 ± 1.0	-0.45 ± 0.11
J0135+3631	01 35 25.2841	+36 31 37.572	82.0 ± 4.1	48.9 ± 1.5	27.0 ± 4.0	16.9 ± 0.9	-0.48 ± 0.12
J0136+3905	01 36 32.5946	+39 05 59.198	84.0 ± 4.0	60.6 ± 1.9	49.0 ± 5.0	42.4 ± 2.1	-0.17 ± 0.09
J0136+3545	01 36 43.6844	+35 45 31.174	115.0 ± 4.0	53.7 ± 1.7	33.0 ± 4.0	22.4 ± 1.1	-0.39 ± 0.10
J0136+4401	01 36 47.3138	+44 01 10.216	97.0 ± 3.9	56.2 ± 1.7	36.0 ± 5.0	21.7 ± 1.1	-0.36 ± 0.11
J0143+3705	01 43 02.5385	+37 05 16.090	80.0 ± 3.7	47.5 ± 1.5	46.0 ± 5.0	40.5 ± 2.0	-0.03 ± 0.09
J0151+4332	01 51 18.3781	+43 32 00.547	42.0 ± 4.1	35.0 ± 1.1	34.0 ± 5.0	26.4 ± 1.3	-0.02 ± 0.12
J0151+4417	01 51 20.8793	+44 17 35.896	74.0 ± 4.3	57.9 ± 2.2	44.0 ± 5.0	62.7 ± 3.1	-0.22 ± 0.10
J0153+4146	01 53 42.2369	+41 46 45.167	250.0 ± 3.7	88.6 ± 3.2	49.0 ± 5.0	33.3 ± 1.7	-0.48 ± 0.09
J0154+3558	01 54 45.4614	+35 58 04.630	< 18.0	30.0 ± 1.3	34.0 ± 5.0	20.9 ± 1.1	+0.10 ± 0.12
J0154+4433	01 54 54.4682	+44 33 37.969	< 18.0	37.7 ± 1.2	40.0 ± 5.0	38.9 ± 2.0	+0.05 ± 0.10
J0156+4459	01 56 28.5219	+44 59 56.484	159.0 ± 4.3	83.5 ± 3.0	46.0 ± 5.0	28.6 ± 1.4	-0.48 ± 0.09
J0157+4120	01 57 05.0059	+41 20 30.634	72.0 ± 3.7	46.5 ± 1.4	30.0 ± 4.0	23.2 ± 1.2	-0.35 ± 0.11
J0159+4059	01 59 14.4871	+40 59 40.403	52.0 ± 3.9	35.2 ± 1.1	33.0 ± 5.0	27.0 ± 1.4	-0.05 ± 0.12
J0159+4144	01 59 49.3346	+41 44 31.946	< 18.0	51.6 ± 1.9	33.0 ± 5.0	21.9 ± 1.1	-0.36 ± 0.13
J1702+5511	17 02 34.5570	+55 11 12.432	< 18.0	38.5 ± 1.2	34.0 ± 4.0	29.5 ± 1.5	-0.10 ± 0.10
J1702+5812	17 02 41.3754	+58 13 10.082	20.0 ± 4.2	40.8 ± 1.3	38.0 ± 5.0	27.7 ± 1.4	-0.06 ± 0.11
J1711+5803	17 11 38.1200	+58 03 28.200	80.0 ± 3.7	44.1 ± 1.8	25.0 ± 4.0		-0.46 ± 0.13
J1715+5724	17 15 22.9752	+57 24 40.312	81.0 ± 3.8	57.6 ± 1.8	35.0 ± 4.0	26.0 ± 1.3	-0.40 ± 0.10
J1720+5541	17 20 09.7600	+55 41 27.300	31.0 ± 3.5	7.9 ± 0.5	27.0 ± 5.0		+0.99 ± 0.16
J1720+5926	17 21 00.6468	+59 26 49.442	137.0 ± 3.7	65.0 ± 2.4	41.0 ± 5.0	22.9 ± 1.2	-0.37 ± 0.10
J1726+5826	17 26 35.1256	+58 26 47.698	43.0 ± 3.8	48.7 ± 1.5	31.0 ± 4.0	38.4 ± 1.9	-0.36 ± 0.11
J1728+5532	17 28 11.6405	+55 32 30.471	21.0 ± 3.6	28.9 ± 1.0	42.0 ± 5.0	33.3 ± 1.7	+0.30 ± 0.10
J1735+5650	17 35 13.7695	+56 50 21.803	86.0 ± 4.1	49.2 ± 2.1	31.0 ± 4.0	33.2 ± 1.7	-0.37 ± 0.11
J1746+5659	17 46 59.5600	+56 59 11.000	< 18.0	12.6 ± 0.6	27.0 ± 4.0		+0.61 ± 0.13
J1747+5902	17 47 33.9322	+59 02 47.967	47.0 ± 4.3	31.9 ± 1.4	25.0 ± 4.0	21.7 ± 1.1	-0.20 ± 0.13
J1756+5918	17 56 11.8346	+59 18 56.824	26.0 ± 4.0	30.1 ± 1.3	25.0 ± 4.0	19.9 ± 1.0	-0.15 ± 0.13
J1756+5806	17 56 29.1449	+58 06 58.236	100.0 ± 3.8	51.6 ± 2.0	38.0 ± 5.0	24.3 ± 1.2	-0.25 ± 0.11
J1758+5951	17 58 12.1100	+59 51 56.100	< 18.0	31.4 ± 1.0	26.0 ± 4.0		-0.15 ± 0.13

**Table A2.** The flux densities and spectral indices from the VLA observations at 4.86, 8.46, 14.94 and 22.46 GHz. The 325 MHz and 1.4 GHz flux densities are taken from WENSS and NVSS, respectively. The letters after the source name correspond to the date of the VLA observations, where *a* is 2000 June 25, *b* is 2001 September 14 and *c* is 2001 December 31.

Source Name	$S_{4.86}$ (mJy)	$S_{8.46}$ (mJy)	$S_{14.94}$ (mJy)	$S_{22.46}$ (mJy)	$\alpha_{0.325}^{1.4}$	$\alpha_{1.4}^{4.86}$	$\alpha_{4.86}^{8.46}$	$\alpha_{8.46}^{14.94}$	$\alpha_{14.94}^{22.46}$
J0103+4322 <sup>b</sup>	41.3 ± 2.1	45.1 ± 2.3	48.1 ± 2.6	39.7 ± 3.2	-0.01 ± 0.07	-0.06 ± 0.06	+0.16 ± 0.13	+0.11 ± 0.13	-0.47 ± 0.24
J0104+3840 <sup>b</sup>	23.7 ± 1.2	17.8 ± 0.9	14.5 ± 1.1	15.4 ± 2.4	> +0.56	-0.43 ± 0.06	-0.52 ± 0.14	-0.36 ± 0.16	+0.15 ± 0.42
J0105+3533 <sup>a</sup>	19.0 ± 1.0	12.8 ± 0.7	9.3 ± 0.7	8.5 ± 0.7	-0.29 ± 0.04	-0.66 ± 0.06	-0.71 ± 0.14	-0.56 ± 0.16	-0.22 ± 0.27
J0106+3522 <sup>a</sup>	22.7 ± 1.2	16.5 ± 0.9	12.3 ± 0.8	8.2 ± 0.7	> +0.56	-0.47 ± 0.06	-0.58 ± 0.14	-0.52 ± 0.14	-0.99 ± 0.25
J0106+4422 <sup>b</sup>	43.4 ± 2.2	26.9 ± 1.4	18.0 ± 1.2	19.1 ± 2.2	+0.54 ± 0.09	-0.42 ± 0.06	-0.86 ± 0.13	-0.71 ± 0.15	+0.15 ± 0.32
J0116+4420 <sup>b</sup>	38.4 ± 1.9	32.2 ± 1.6	21.6 ± 1.4	15.6 ± 2.0	-0.61 ± 0.04	-0.48 ± 0.06	-0.32 ± 0.13	-0.70 ± 0.14	-0.80 ± 0.35
J0119+4054 <sup>b</sup>	56.6 ± 2.9	54.4 ± 2.7	47.6 ± 2.5	43.8 ± 2.8	-0.09 ± 0.05	-0.23 ± 0.06	-0.07 ± 0.13	-0.23 ± 0.13	-0.20 ± 0.20
J0123+3843 <sup>a</sup>	35.2 ± 1.8	33.0 ± 1.7	21.4 ± 1.2	15.0 ± 0.9	> +0.02	+0.52 ± 0.06	-0.12 ± 0.13	-0.76 ± 0.13	-0.87 ± 0.20
J0123+3806 <sup>b</sup>	86.1 ± 4.3	84.2 ± 4.2	64.0 ± 3.3	47.5 ± 2.9	-0.15 ± 0.07	+0.68 ± 0.06	-0.04 ± 0.13	-0.48 ± 0.13	-0.73 ± 0.19
J0123+3641 <sup>a</sup>	42.4 ± 2.2	29.7 ± 1.5	19.6 ± 1.1	12.3 ± 0.8	+0.30 ± 0.08	-0.19 ± 0.06	-0.64 ± 0.13	-0.73 ± 0.13	-1.14 ± 0.21
J0128+4003 <sup>b</sup>	40.0 ± 2.0	31.5 ± 1.6	20.5 ± 1.3	< 3.0	-0.49 ± 0.04	-0.27 ± 0.06	-0.43 ± 0.13	-0.76 ± 0.14	< -4.70
J0128+4439 <sup>b</sup>	38.7 ± 2.0	35.2 ± 1.8	31.0 ± 1.8	26.6 ± 2.4	> +0.49	+0.04 ± 0.06	-0.17 ± 0.13	-0.22 ± 0.13	-0.37 ± 0.26
J0129+4044 <sup>b</sup>	35.8 ± 1.8	30.5 ± 1.6	32.9 ± 1.9	31.3 ± 2.6	-0.65 ± 0.04	-0.51 ± 0.06	-0.29 ± 0.13	+0.12 ± 0.13	-0.12 ± 0.25
J0131+4428 <sup>b</sup>	30.0 ± 1.5	36.5 ± 1.9	38.4 ± 2.2	33.9 ± 2.7	-0.45 ± 0.05	-0.10 ± 0.06	+0.35 ± 0.14	+0.09 ± 0.13	-0.31 ± 0.24
J0132+4345 <sup>b</sup>	43.7 ± 2.2	37.9 ± 1.9	35.3 ± 2.0	33.1 ± 2.8	-0.40 ± 0.04	-0.18 ± 0.06	-0.26 ± 0.13	-0.13 ± 0.13	-0.16 ± 0.25
J0134+4018 <sup>b</sup>	30.5 ± 1.6	22.8 ± 1.2	18.0 ± 0.9	< 3.0	-0.49 ± 0.04	-0.41 ± 0.06	-0.52 ± 0.14	-0.42 ± 0.13	< -4.39
J0135+3631 <sup>b</sup>	23.3 ± 1.2	17.2 ± 0.9	13.7 ± 0.7	< 3.0	-0.35 ± 0.05	-0.60 ± 0.06	-0.55 ± 0.14	-0.40 ± 0.13	< -3.73
J0136+3905 <sup>a</sup>	52.0 ± 2.6	48.3 ± 2.4	46.8 ± 2.4	44.5 ± 2.3	-0.22 ± 0.05	-0.12 ± 0.06	-0.13 ± 0.13	-0.06 ± 0.13	-0.12 ± 0.18
J0136+3545 <sup>b</sup>	22.5 ± 1.2	16.8 ± 0.9	13.9 ± 1.4	< 3.0	-0.52 ± 0.04	-0.70 ± 0.06	-0.53 ± 0.14	-0.33 ± 0.20	< -3.76
J0136+4401 <sup>b</sup>	29.3 ± 1.5	18.7 ± 1.0	14.7 ± 1.2	< 3.0	-0.37 ± 0.04	-0.52 ± 0.06	-0.81 ± 0.14	-0.42 ± 0.17	< -3.90
J0143+3705 <sup>b</sup>	53.6 ± 2.7	46.2 ± 2.3	29.2 ± 1.9	20.5 ± 3.0	-0.36 ± 0.05	+0.10 ± 0.06	-0.27 ± 0.13	-0.81 ± 0.14	-0.87 ± 0.39
J0151+4332 <sup>b</sup>	33.9 ± 1.7	34.4 ± 1.8	31.7 ± 1.8	28.5 ± 2.4	-0.12 ± 0.08	-0.03 ± 0.06	+0.03 ± 0.14	-0.14 ± 0.13	-0.26 ± 0.25
J0151+4417 <sup>b</sup>	85.4 ± 4.3	72.3 ± 3.6	48.3 ± 2.6	34.9 ± 2.9	-0.17 ± 0.05	+0.31 ± 0.06	-0.30 ± 0.13	-0.71 ± 0.13	-0.80 ± 0.24
J0153+4146 <sup>a</sup>	25.6 ± 1.3	32.5 ± 1.7	35.5 ± 1.9	34.8 ± 1.9	-0.71 ± 0.04	-1.00 ± 0.06	+0.43 ± 0.14	+0.16 ± 0.13	-0.05 ± 0.18
J0154+3558 <sup>b</sup>	30.3 ± 1.6	19.8 ± 1.0	12.8 ± 1.3	< 3.0	> +0.35	+0.01 ± 0.06	-0.77 ± 0.14	-0.77 ± 0.20	< -3.56
J0154+4433 <sup>a</sup>	48.6 ± 2.4	46.2 ± 2.3	44.9 ± 2.3	37.4 ± 2.0	> +0.51	+0.20 ± 0.06	-0.09 ± 0.13	-0.05 ± 0.13	-0.45 ± 0.18
J0156+4459 <sup>b</sup>	40.5 ± 2.1	27.2 ± 1.4	17.5 ± 1.2	16.3 ± 2.0	-0.44 ± 0.04	-0.58 ± 0.06	-0.72 ± 0.14	-0.78 ± 0.15	-0.17 ± 0.35
J0157+4120 <sup>a</sup>	29.8 ± 1.5	22.8 ± 1.1	19.3 ± 1.1	15.8 ± 1.0	-0.30 ± 0.05	-0.36 ± 0.06	-0.48 ± 0.14	-0.29 ± 0.13	-0.49 ± 0.21
J0159+4059 <sup>b</sup>	38.2 ± 1.9	30.1 ± 1.5	21.8 ± 1.4	15.8 ± 1.7	-0.27 ± 0.05	+0.07 ± 0.06	-0.43 ± 0.14	-0.57 ± 0.14	-0.79 ± 0.30
J0159+4144 <sup>b</sup>	30.1 ± 1.5	23.0 ± 1.2	12.6 ± 1.2	< 3.0	> +0.72	-0.43 ± 0.06	-0.49 ± 0.14	-1.06 ± 0.18	< -3.52
J1702+5511 <sup>c</sup>	25.3 ± 1.4	22.5 ± 1.1	17.6 ± 1.0	15.1 ± 0.8	> +0.52	-0.34 ± 0.06	-0.21 ± 0.14	-0.43 ± 0.13	-0.38 ± 0.19
J1702+5812 <sup>c</sup>	34.2 ± 1.7	28.0 ± 1.4	20.8 ± 1.1	16.2 ± 0.9	+0.49 ± 0.15	-0.14 ± 0.06	-0.36 ± 0.13	-0.52 ± 0.13	-0.61 ± 0.19
J1711+5803 <sup>a</sup>	21.2 ± 1.1	12.4 ± 0.7	6.4 ± 0.6	< 3.0	-0.41 ± 0.05	-0.59 ± 0.06	-0.97 ± 0.15	-1.16 ± 0.20	< -1.86
J1715+5724 <sup>c</sup>	28.6 ± 1.5	20.4 ± 1.0	14.6 ± 0.8	11.7 ± 0.7	-0.23 ± 0.05	-0.56 ± 0.06	-0.61 ± 0.13	-0.59 ± 0.13	-0.54 ± 0.20
J1720+5541 <sup>a</sup>	25.8 ± 1.3	22.4 ± 1.2	13.6 ± 0.9	10.8 ± 0.8	-0.94 ± 0.08	+0.95 ± 0.06	-0.25 ± 0.14	-0.88 ± 0.15	-0.57 ± 0.25
J1720+5926 <sup>c</sup>	32.7 ± 1.7	21.8 ± 1.1	17.0 ± 1.0	13.6 ± 0.8	-0.51 ± 0.04	-0.55 ± 0.06	-0.73 ± 0.13	-0.44 ± 0.13	-0.55 ± 0.19
J1726+5826 <sup>a</sup>	43.6 ± 2.2	38.9 ± 2.0	35.9 ± 1.9	33.4 ± 1.8	+0.09 ± 0.07	-0.09 ± 0.06	-0.21 ± 0.13	-0.14 ± 0.13	-0.18 ± 0.18
J1728+5532 <sup>c</sup>	42.7 ± 2.2	40.7 ± 2.0	38.1 ± 2.0	33.5 ± 1.7	+0.22 ± 0.12	+0.31 ± 0.06	-0.09 ± 0.13	-0.12 ± 0.13	-0.32 ± 0.18
J1735+5650 <sup>c</sup>	39.6 ± 2.0	40.3 ± 2.0	39.4 ± 2.0	35.6 ± 1.8	-0.38 ± 0.05	-0.17 ± 0.06	+0.03 ± 0.13	-0.04 ± 0.13	-0.25 ± 0.18
J1746+5659 <sup>c</sup>	28.6 ± 1.5	34.5 ± 1.7	23.0 ± 1.2	14.7 ± 0.8	> -0.24	+0.66 ± 0.06	+0.34 ± 0.13	-0.71 ± 0.13	-1.10 ± 0.19
J1747+5902 <sup>a</sup>	33.3 ± 1.7	36.3 ± 1.8	45.2 ± 2.3	43.2 ± 2.2	-0.27 ± 0.07	+0.03 ± 0.06	+0.16 ± 0.14	+0.39 ± 0.13	-0.11 ± 0.18
J1756+5918 <sup>c</sup>	25.3 ± 1.3	23.7 ± 1.2	24.2 ± 1.3	24.0 ± 1.2	+0.10 ± 0.11	-0.14 ± 0.06	-0.12 ± 0.14	+0.04 ± 0.13	-0.02 ± 0.18
J1756+5806 <sup>c</sup>	31.3 ± 1.6	24.4 ± 1.2	18.3 ± 1.0	15.7 ± 0.9	-0.45 ± 0.04	-0.40 ± 0.06	-0.45 ± 0.13	-0.51 ± 0.13	-0.38 ± 0.19
J1758+5951 <sup>a</sup>	53.5 ± 2.7	65.8 ± 3.3	36.3 ± 1.9	26.5 ± 1.5	> +0.38	+0.43 ± 0.06	+0.37 ± 0.13	-1.05 ± 0.13	-0.77 ± 0.19

## B Source code

Relevant source code is listed here.

### B.1 Least Squares fitting algorithm

The following script is used to fit all spectral types using an LSq algorithm.

Scripts/Thesis\_Sv-omnifit-LSq\_v2-1.py

```
1 #!/usr/bin/env python
2 from __future__ import division
3 import numpy as np
4 import scipy as sp
5 import scipy.interpolate as spi
6 import scipy.optimize as spo
7 import scipy.stats as sps
8 from matplotlib.pyplot import figure, show, rc, rcParams
9 #####THIS SCRIPT CAN RUN ALL THREE FITS, BUT UNCERTAINTIES HAVE BEEN REPRESSED BY SOMEWHAT
10 ARGUABLE MEANS. ONLY THE SYNCH PLUS POWER LAW SPECTRUM'S UNCERTAINTIES STAND IN MY
11 PATH TOWARDS POWER, INFINITE POWER!
12 #####This script utilizes a least squares algorithm to obtain the best fit.
13 #####Episode I: Reading Data (or: The Phantom Data).
14 #####
15 print(">>> Episode I: The Phantom Data")
16 #####Reading relevant data from (mono-axial) datafile. If uncertainties were unknown, they
17 have been set to 1.0 to prevent code collapse due to divide by zero errors.
18 fulllist = np.genfromtxt("Janskylist.dat", dtype=None, usecols=[0], comments="#")
19 errlist = np.genfromtxt("Janskylist.dat", dtype=None, usecols=[1], comments="#")
20 #####Splitting the superarray into useful, seperate arrays with relevant data.
21 names, S003, S014, S048, S084, S149, S224 = np.split(fulllist, 7)
22 junkerr, errS003, errS014, errS048, errS084, errS149, errS224 = np.split(errlist, 7)
23 names = names.astype(str).flatten()
24 S003 = np.asarray(S003.astype(float).flatten())
25 S014 = np.asarray(S014.astype(float).flatten())
26 S048 = np.asarray(S048.astype(float).flatten())
27 S084 = np.asarray(S084.astype(float).flatten())
28 S149 = np.asarray(S149.astype(float).flatten())
29 S224 = np.asarray(S224.astype(float).flatten())
30 #####Printing all data in a clear-cut table.
31 print("\033[1m", "index * name * 0.325 * 1.4 * 4.85 * 8.46 * 14.94 * 22.46", "\033[0m",
32 <- Frequency bands")
33 for i in range(len(names)):
34     print(i, ":", "\033[1m", names[i], "\033[0m", "*", S003[i], "*", S014[i], "*", S048[i], "*",
35     S084[i], "*", S149[i], "*", S224[i])
36 n = int(input(" Pick GPS source: "))
37 #####Creating the to be used data arrays.
38 v = [0.325, 1.4, 4.85, 8.46, 14.94, 22.46]
39 S = [S003[n], S014[n], S048[n], S084[n], S149[n], S224[n]]
40 Serr = [errS003[n], errS014[n], errS048[n], errS084[n], errS149[n], errS224[n]]
41 print("Source designation:", names[n])
42 print("Frequency bands:", np.asarray(v))
43 print("Flux densities:", np.asarray(S))
44 print("Flux uncertainties:", np.asarray(Serr))
45 #####Plotting the recorded data points for visualisation purposes.
46 fig = figure(figsize=(10,10))
47 frame = fig.add_subplot(1,1,1, xscale="log", yscale="log")
48 frame.errorbar(v, S, yerr=Serr, ecolor="k", marker="*", mfc="y", ms=10, ls="", label="Data from
49 {}".format(names[n]))
50 frame.set_xlabel('Frequency v (GHz)')
51 frame.set_ylabel('Flux density S (mJy)')
52 frame.legend()
```

```

53 show()
54
55 ###While-loop to allow user-friendly cancelling of code.
56 while True:
57     print("\033[1m","Do you want to continue?","\033[0m")
58     answer = input(" :")
59     if answer.lower().startswith("y"):
60         print("Roger")
61         break
62     elif answer.lower().startswith("n"):
63         print("OKDOEP")
64         exit()
65
66 ###While-loop to allow user-friendly usage of power law or not, with q the model index,
67     ndim the amount of free parameters, and os the offset within the naming list for
68     easier printing of results. For q = 3, the spectrum shows signs of ageing – a final
69     point that is offset from the power law-like behaviour of its predecessors. Fitting
70     an ageing component is extremely difficult, so we instead ignore the final point and
71     plot a power law through the others as spectral ageing is not of interest to our
72     research.
73 while True:
74     print("\033[1m","What kind of fit would you like to use?","\033[0m")
75     print("0 = power law, 1 = synchrotron spectrum, 2 = synchrotron + power law, 3 =
76     ageing power law spectrum, 4 = synchrotron + underdefined power law (i.e. only
77     through first two points), 5 = double power law.")
78     answer = input(" :")
79     if answer.lower().startswith("0"):
80         q = 0 ; ndim = 2 ; os=4
81         print("Roger: power law online")
82         break
83     elif answer.lower().startswith("1"):
84         q = 1 ; ndim = 4 ; os=0
85         print("Roger: synch fit online")
86         break
87     elif answer.lower().startswith("2"):
88         q = 2 ; ndim = 6 ; os=0
89         print("Roger: synch fit and power law online")
90         break
91     elif answer.lower().startswith("3"):
92         q = 3 ; ndim = 2 ; os=4
93         print("Roger: ageing fit online. Final datapoint will be ignored")
94         break
95     elif answer.lower().startswith("4"):
96         q = 4 ; ndim = 5 ; os=0
97         print("Roger: synch fit and undef power law online. A fudgy powerlaw will be
98         drawn.")
99         break
100    elif answer.lower().startswith("5"):
101        q = 5 ; ndim = 4 ; os=4
102        print("Roger: double power law online.")
103        break
104    modelname = ["power law","synchrotron","synchrotron/power law","ageing spectrum",
105        "synchrotron/undef power law","double power law"]
106    print(" Number of free parameters:" ,ndim)
107
108    #####Episode II: Creating fitting functions and derivatives (or: Attack of the Functions
109    ). #####
110    print(">>> Episode II: Attack of the Functions")
111
112    ###Defining the fits.
113    def Sfitpl(v,A,m):
114        f1 = A*v**m                ###First power law
115        return f1
116
117    def Sfitplpl(v,A1,m1,A2,m2):
118        f1 = A1*v**m1                ###First power law
119        f2 = A2*v**m2                ###First power law
120        return f1+f2

```

```

112 def Sfitffa(v,A,m,t):
113     f1 = A*v**m                                     ###First power law
114     f2 = np.e**(-t*v**(-2.1))                     ###Exponential fall-off
115     return f1*f2
116
117 def Sfitsyn(v,S01,v01,k,l):
118     f1 = S01*(v/v01)**k                             ###First power law
119     f2 = (1-np.e**(-(v/v01)**(1-k)))/(1-np.e**(-1))  ###Second power law (exponential
120     fall-off)
121     return f1*f2
122
123 def Sfitplsyn(v,S01,v01,k,l,A,m):
124     f1 = S01*(v/v01)**k                             ###First power law
125     f2 = (1-np.e**(-(v/v01)**(1-k)))/(1-np.e**(-1))  ###Second power law (exponential
126     fall-off)
127     f3 = A*v**m                                     ###Third power law (only when two
128     turnovers are present) Wherein A=S02/(v02^m)
129     return f1*f2 + f3
130
131 def derSfitsyn(v,S01,v01,k,l):                     ###Derivative of synchrotron fit
132     E = np.e**(-(v/v01)**(1-k))
133     C = S01/(v01*(1-1/np.e))
134     f1 = C*k*((v/v01)**(k-1))*(1-E)
135     f2 = C*(1-k)*((v/v01)**(1-1))*E
136     return f1 + f2
137
138 def derSfitplsyn(v,S01,v01,k,l,A,m):              ###Derivative of power law/
139     synchrotron combined fit
140     E = np.e**(-(v/v01)**(1-k))
141     C = S01/(v01*(1-1/np.e))
142     f1 = C*k*((v/v01)**(k-1))*(1-E)
143     f2 = C*(1-k)*((v/v01)**(1-1))*E
144     f3 = A*m*v**(m-1)
145     return f1 + f2 + f3
146
147 ###Array spanning the fitting space.
148 vfit = np.linspace(0.3,23,1000)
149
150 ###List that names the coefficients for the synchfit.
151 coeflist = ["S_{01}", "v_{01}", "k", "l", "A1", "m1", "A2", "m2"]
152 ###With S_{01} the scale flux of the synchrotron model, v_{01} the scale frequency of
153     the synchrotron model, k the optically thick spectral index (before the peak), l the
154     optically thin spectral index (after the peak), A = S_{02}/(v_{02}^m) the fudged
155     constant for the power law, combining the power law scale flux and scale frequency (
156     S_{02} and v_{02}, respectively) with the power law spectral index k=l -> m (
157     otherwise known as Jesper's Fudge Factor), and m the power law spectral index for a
158     spectrum wherein k=l.
159
160 #####Episode III: Computing uncertainties in fits. (Revenge of the Uncertainties).
161     #####
162 print(">>> Episode III: Revenge of the Uncertainties")
163
164 def errSfitpl(v,A,m,cov):                          ###Finding uncertainties function in power law
165     .
166     dsdA = v**m
167     dsdm = A*(v**m)*np.log(v)
168     dsdx = np.asarray([dsdA, dsdm])
169     Svar = 0
170     for i in range(ndim):
171         for j in range(ndim):
172             Svar = Svar + dsdx[i]*dsdx[j]*cov[i,j]
173     return np.sqrt(Svar)
174
175 def errSfitplpl(v,A1,m1,A2,m2,cov):                ###Finding uncertainties function in
176     power law.
177     dsdA1 = v**m1
178     dsdm1 = A1*(v**m1)*np.log(v)
179     dsdA2 = v**m2
180     dsdm2 = A2*(v**m2)*np.log(v)

```

```

169 dsdx = np. asarray ([dsdA1, dsdm1, dsdA2, dsdm2])
170 Svar = 0
171 for i in range(ndim):
172     for j in range(ndim):
173         Svar = Svar + dsdx[i]*dsdx[j]*cov[i,j]
174 return np. sqrt(Svar)
175
176 def errSfitffa(v,A,m,t,cov):          ###Finding uncertainties function in power law
177 .
178 dsdA = (v**m)*np. e**(-t*v**-2.1)
179 dsdm = A*(v**m)*np. log(v)*np. e**(-t*v**-2.1)
180 dsdt = -A*(v**(m-2.1))*np. e**(-t*v**-2.1)
181 dsdx = np. asarray ([dsdA, dsdm, dsdt])
182 Svar = 0
183 for i in range(ndim):
184     for j in range(ndim):
185         Svar = Svar + dsdx[i]*dsdx[j]*cov[i,j]
186 return np. sqrt(Svar)
187
188 def errSfitsyn(v,S01,v01,k,l,cov):    ###Finding uncertainties function in synch fit
189 .
190 C = S01/(np. e-1)
191 C2 = S01/(1-1/np. e)
192 E = np. e**(-(v/v01)**(1-k))
193 E2 = np. e**(1-((v/v01)**(1-k)))
194 dsdS01 = Sfitsyn(v,S01,v01,k,l)/S01
195 dsdv01 = (-((C2*(1-E)))*k*v*((v/v01)**(k-1)))/(v01**2) - (C2*(1-k)*v*E*((v/v01)**(1-1)))/(v01**2)
196 dsdk = C2*(1-E)*((v/v01)**k)*np. log(v/v01) - C2*((v/v01)**(1))*np. log(v/v01)*E
197 dsdl = C*((v/v01)**1)*np. log(v/v01)*E2
198 dsdx = np. asarray ([dsdS01, dsdv01, dsdk, dsdl])
199 Svar = 0
200 for i in range(ndim):
201     for j in range(ndim):
202         Svar = Svar + dsdx[i]*dsdx[j]*cov[i,j]
203 return np. sqrt(Svar)
204
205 def errSfitplsyn(v,S01,v01,k,l,A,m,cov): ###Finding uncertainties function in synch fit
206 plus power law.
207 C = S01/(np. e-1)
208 C2 = S01/(1-1/np. e)
209 E = np. e**(-(v/v01)**(1-k))
210 E2 = np. e**(1-((v/v01)**(1-k)))
211 dsdS01 = Sfitsyn(v,S01,v01,k,l)/S01
212 dsdv01 = (-((C2*(1-E)))*k*v*((v/v01)**(k-1)))/(v01**2) - (C2*(1-k)*v*E*((v/v01)**(1-1)))/(v01**2)
213 dsdk = C2*(1-E)*((v/v01)**k)*np. log(v/v01) - C2*((v/v01)**(1))*np. log(v/v01)*E
214 dsdl = C*((v/v01)**1)*np. log(v/v01)*E2
215 dsdA = v**m
216 dsdm = A*(v**m)*np. log(v)
217 dsdx = np. asarray ([dsdS01, dsdv01, dsdk, dsdl, dsdA, dsdm])
218 Svar = 0
219 for i in range(ndim):
220     for j in range(ndim):
221         Svar = Svar + dsdx[i]*dsdx[j]*cov[i,j]
222 return np. sqrt(Svar)
223
224 #####Episode IV: Least squares fitting the data (or: A New Fit).
225 #####
226 print(">>> Episode IV: A New Fit")
227
228 ###Creating some new, useful arrays and indices.
229 S2 = [S[1],S[2],S[3],S[4],S[5]]
230 Serr2 = [Serr[1],Serr[2],Serr[3],Serr[4],Serr[5]]
231 v2 = [v[1],v[2],v[3],v[4],v[5]]
232 S3 = [S[0],S[1],S[2],S[3],S[4]]
233 Serr3 = [Serr[0],Serr[1],Serr[2],Serr[3],Serr[4]]
234 v3 = [v[0],v[1],v[2],v[3],v[4]]
235 Smi = S. index(max(S))

```

```

233 Smi2 = S2.index(max(S2))
234 S1 = [S[0],S[1],S[2]]
235 v1 = [v[0],v[1],v[2]]
236 Smi1 = S1.index(max(S1))
237 Sh = [S[3],S[4],S[5]]
238 vh = [v[3],v[4],v[5]]
239 Smih = Sh.index(max(Sh))
240
241 ###Monstrous, hulking, hideous if-loop that creates the fit by first estimating and then
least squaring parameters.
242 if q == 0:
243     S0guess = max(S)
244     v0guess = v[Smi]
245     mguess = (np.log10(S[5])-np.log10(S[0]))/(np.log10(v[5])-np.log10(v[0]))
246     Aguess = S0guess/(v0guess**mguess)
247     estimates = [Aguess,mguess]
248     p, pcov = spo.curve_fit(Sfitpl,v,S,[Aguess,mguess],sigma=Serr,bounds
249     =([0,-10],[500,10]),**{'max_nfev':1000000})
250     fitfun = Sfitpl(vfit,p[0],p[1])
251     fitmax = fitfun + errSfitpl(vfit,p[0],p[1],pcov)
252     fitmin = fitfun - errSfitpl(vfit,p[0],p[1],pcov)
253 elif q == 1:
254     S0guess = max(S)
255     v0guess = v[Smi]
256     while True:
257         print("\033[1m","Do you want to constrain k narrowly to 2.3-2.5?","\033[0m")
258         answer = input(" :")
259         if answer.lower().startswith("y"):
260             klow = 2.3
261             kguess = 2.4
262             print("Roger: k constrained to 2.3-2.5")
263             break
264         if answer.lower().startswith("n"):
265             klow = 0.0
266             if S[1] > S[0]:
267                 kguess = (np.log10(S[1])-np.log10(S[0]))/(np.log10(v[1])-np.log10(v[0]))
268             else:
269                 if S[2] > S[0]:
270                     kguess = (np.log10(S[2])-np.log10(S[0]))/(np.log10(v[2])-np.log10(v
271 [0]))
272                 else:
273                     kguess = 2.0
274             print("Roger: k not narrowly constrained.")
275             break
276         if S[5] < S[4]:
277             lguess = (np.log10(S[5])-np.log10(S[4]))/(np.log10(v[5])-np.log10(v[4]))
278         else:
279             if S[5] < S[3]:
280                 lguess = (np.log10(S[5])-np.log10(S[3]))/(np.log10(v[5])-np.log10(v[3]))
281             else:
282                 lguess = -1.0
283     estimates = [S0guess,v0guess,kguess,lguess]
284     p, pcov = spo.curve_fit(Sfitsyn,v,S,[S0guess,v0guess,kguess,lguess],sigma=Serr,
285     bounds=([0,0.3,klow,-5.0],[500,23,2.5,0]),**{'max_nfev':1000000})
286     fitfun = Sfitsyn(vfit,p[0],p[1],p[2],p[3])
287     fitmax = fitfun + errSfitsyn(vfit,p[0],p[1],p[2],p[3],pcov)
288     fitmin = fitfun - errSfitsyn(vfit,p[0],p[1],p[2],p[3],pcov)
289 elif q == 2:
290     S0guess = max(S2)
291     v0guess = v2[Smi2]
292     kguess = (np.log10(S2[Smi2])-np.log10(S2[Smi2-1]))/(np.log10(v2[Smi2])-np.log10(v2[
293 Smi2-1]))
294     if S[5] < S[4]:
295         lguess = (np.log10(S[5])-np.log10(S[4]))/(np.log10(v[5])-np.log10(v[4]))
296     else:
297         if S[5] < S[3]:
298             lguess = (np.log10(S[5])-np.log10(S[3]))/(np.log10(v[5])-np.log10(v[3]))
299         else:
300             lguess = -1.0
301     Aguess = max(S)

```

```

298 if S[1] < S[0]:
299     mguess = (np.log10(S[1])-np.log10(S[0]))/(np.log10(v[1])-np.log10(v[0]))
300 else:
301     if S[2] < S[0]:
302         mguess = (np.log10(S[2])-np.log10(S[0]))/(np.log10(v[2])-np.log10(v[0]))
303     else:
304         mguess = -1.0
305     estimates = [S0guess, v0guess, kguess, lguess, Aguess, mguess]
306     p, pcov = spo.curve_fit(Sfitplsyn, v, S, [S0guess, v0guess, kguess, lguess, Aguess, mguess],
307                             sigma=Serr, bounds=([0, 0.3, 0, -5.0, 0, -10], [500, 23, 2.5, 0, 500, 10]), **{'max_nfev':
308                             :1000000})
309     fitfun = Sfitplsyn(vfit, p[0], p[1], p[2], p[3], p[4], p[5])
310     fitmax = fitfun + errSfitplsyn(vfit, p[0], p[1], p[2], p[3], p[4], p[5], pcov)
311     fitmin = fitfun - errSfitplsyn(vfit, p[0], p[1], p[2], p[3], p[4], p[5], pcov)
312 elif q == 3:
313     S0guess = max(S)
314     v0guess = v[Smi]
315     mguess = (np.log10(S[4])-np.log10(S[0]))/(np.log10(v[4])-np.log10(v[0]))
316     Aguess = S0guess/(v0guess**mguess)
317     estimates = [Aguess, mguess]
318     p, pcov = spo.curve_fit(Sfitpl, v3, S3, [Aguess, mguess], sigma=Serr3, bounds
319     =([0, -10], [500, 0]), **{'max_nfev':1000000})
320     fitfun = Sfitpl(vfit, p[0], p[1])
321     fitmax = fitfun + errSfitpl(vfit, p[0], p[1], pcov)
322     fitmin = fitfun - errSfitpl(vfit, p[0], p[1], pcov)
323 elif q == 4:
324     S0guess = max(S2)
325     v0guess = v[Smi2]
326     if S2[1] > S2[0]:
327         kguess = (np.log10(S2[1])-np.log10(S2[0]))/(np.log10(v2[1])-np.log10(v2[0]))
328     else:
329         if S2[2] > S2[0]:
330             kguess = (np.log10(S2[2])-np.log10(S2[0]))/(np.log10(v2[2])-np.log10(v2[0]))
331         else:
332             kguess = 2.0
333     if S2[4] < S2[3]:
334         lguess = (np.log10(S2[4])-np.log10(S2[3]))/(np.log10(v2[4])-np.log10(v2[3]))
335     else:
336         if S2[4] < S2[2]:
337             lguess = (np.log10(S2[4])-np.log10(S2[2]))/(np.log10(v2[4])-np.log10(v2[2]))
338         else:
339             lguess = -1.0
340     M = (np.log10(0.5*S[1])-np.log10(S[0]))/(np.log10(v[1])-np.log10(v[0]))
341     Aguess = S[0]/(v[0]**M)
342     estimates = [S0guess, v0guess, kguess, lguess, Aguess, M]
343     def Sfitplsyn2(v, S01, v01, k, l, A):
344         f1 = S01*(v/v01)**k                                     ###First power law
345         f2 = (1-np.e**(-(v/v01)**(1-k)))/(1-np.e**(-1))      ###Second power law (
346         exponential fall-off)
347         f3 = A*v**M                                           ###Third power law (only when two
348         turnovers are present) Wherein A=S02/(v02^m)
349         return f1*f2 + f3
350     p, pcov = spo.curve_fit(Sfitplsyn2, v, S, [S0guess, v0guess, kguess, lguess, Aguess], sigma=
351     Serr, bounds=([0, 0.3, 0, -5.0, 0], [500, 50, 2.5, 0, 500]), **{'max_nfev':1000000})
352     fitfun = Sfitplsyn2(vfit, p[0], p[1], p[2], p[3], p[4], M)
353     fitmax = fitfun + errSfitplsyn2(vfit, p[0], p[1], p[2], p[3], p[4], M, pcov)
354     fitmin = fitfun - errSfitplsyn2(vfit, p[0], p[1], p[2], p[3], p[4], M, pcov)
355 if q == 5:
356     S01guess = max(S1)
357     v01guess = v1[Smi1]
358     S02guess = max(S2)
359     v02guess = v2[Smi2]
360     m1guess = (np.log10(0.5*S[2])-np.log10(S[0]))/(np.log10(v[1])-np.log10(v[0]))
361     A1guess = S01guess/(v01guess**m1guess)
362     m2guess = (np.log10(S[5])-np.log10(0.5*S[3]))/(np.log10(v[5])-np.log10(v[3]))
363     A2guess = S02guess/(v02guess**m2guess)
364     estimates = [A1guess, m1guess, A2guess, m2guess]
365     p, pcov = spo.curve_fit(Sfitplpl, v, S, [A1guess, m1guess, A2guess, m2guess], sigma=Serr,
366     bounds=([0, -10, 0, -10], [500, 10, 500, 10]), **{'max_nfev':1000000})
367     fitfun = Sfitplpl(vfit, p[0], p[1], p[2], p[3])

```

```

361 fitmax = fitfun + errSfitplpl(vfit,p[0],p[1],p[2],p[3],pcov)
362 fitmin = fitfun - errSfitplpl(vfit,p[0],p[1],p[2],p[3],pcov)
363
364
365 ####Printing loop outputs, i.e. values of free parameters by least squares fitting.
366 print(" Parameter estimates are:",estimates)
367 print("\033[1m"+ " >>> Using scipy.optimize.curve_fit on {} model:".format(modelname[q
368 ]),"\033[0m")
369 for i in range(ndim):
370     print(" Coefficient",coeflist[i+os],"equals",p[i])
371     print(" Uncertainty of coefficient",coeflist[i+os],"equals",np.sqrt(pcov[i,i]))
372 if q == 4:
373     print(" Coefficient m (fixed) equals",M)
374
375 #####Episode V: Finding the maximum of the fit (or: The Maximum Strikes Back).
376 #####
377 print(">>> Episode V: The Maximum Strikes Back")
378
379 ####If loop that bisection the function if it has one or more peaks.
380 if q == 0 or q == 3 or q == 5:
381     vmax = 0
382     Smax = 0
383 elif q == 1 or q == 2 or q == 4:
384     if q == 2 or q == 4:
385         if S[2] > S[1]:
386             inc = (v[2]+v[1])/2
387         else:
388             if S[3] > S[2]:
389                 inc = (v[3]+v[2])/2
390             else:
391                 inc = v[3]
392     def bisector(v):
393         #####Function of one variable used by the
394         bisection algorithm.
395         return derSfitplsyn(v,p[0],p[1],p[2],p[3],p[4],M)
396 elif q == 1:
397     inc = v[0]
398     def bisector(v):
399         #####Function of one variable used by the
400         bisection algorithm.
401         return derSfitsyn(v,p[0],p[1],p[2],p[3])
402 vfitbis = np.linspace(inc,23,1000)
403 for i in range(len(vfitbis)-1):
404     if bisector(vfitbis[i+1]) < 0 and bisector(vfitbis[i]) > 0:
405         ma = vfitbis[i+1]
406         mi = vfitbis[i]
407 vmaxerr = (23-inc)/1000 #Stepsize of fitting space.
408 vmax = spo.bisect(bisector,mi,ma,xtol=vmaxerr)
409 if q == 1:
410     Smax = Sfitsyn(vmax,p[0],p[1],p[2],p[3])
411     Smaxerr = errSfitsyn(vmax,p[0],p[1],p[2],p[3],pcov)
412 elif q == 2:
413     Smax = Sfitplsyn(vmax,p[0],p[1],p[2],p[3],p[4],p[5])
414     Smaxerr = errSfitplsyn(vmax,p[0],p[1],p[2],p[3],p[4],p[5],pcov)
415 elif q == 4:
416     Smax = Sfitplsyn(vmax,p[0],p[1],p[2],p[3],p[4],M)
417     Smaxerr = errSfitplsyn(vmax,p[0],p[1],p[2],p[3],p[4],M,pcov)
418
419 #####Printing peak coordinates.
420 print("\033[1m"," >>> Using scipy.optimize.bisect on {} model gives peak coordinates:
421 ".format(modelname[q]),"\033[0m")
422 if q == 0 or q == 3 or q == 5:
423     print(" A peak could not be determined, for a power law model is without one.")
424 else:
425     print(" Fit frequency of peak is:",vmax,"GHz")
426     print(" Uncertainty of frequency peak is:",vmaxerr,"GHz")
427     print(" Fit flux density of peak is:",Smax,"mJy")
428     print(" Uncertainty of flux peak is:",Smaxerr,"mJy")
429
430 #####Episode VI: Creating pdf's from coeff's found in Ep.IV (or: Return of the Pdf).

```

```

#####
426 print(">>> Episode VI: Return of the Pdf")
427
428 def mpdf(m1,m2,s11,s12,s21,s22,x): ###Multivariate pdf maker function.
429     y = sps.multivariate_normal([m1,m2],[[s11,s12],[s21,s22]])
430     return y.pdf(x)
431
432 def mpdfax(m1,m2,s11,s22): ###Multivariate pdf grid maker function.
433     up1 = (m1+3*np.sqrt(s11)) ; up2 = (m2+3*np.sqrt(s22))
434     low1 = (m1-3*np.sqrt(s11)) ; low2 = (m2-3*np.sqrt(s22))
435     x1,x2 = np.mgrid[low1:up1:(up1-low1)/1000, low2:up2:(up2-low2)/1000]
436     pos2 = np.empty(x1.shape + (2,))
437     pos2[:, :, 0] = x1 ; pos2[:, :, 1] = x2
438     return x1,x2,pos2
439
440
441 #####Episode VII: Plotting all results (or: The Pyplot Awakens).
442 #####
442 print(">>> Episode VII: The Pyplot Awakens")
443
444 ###Plotting data, least squares fits and peaks (Episodes I, IV & V).
445 fig1 = figure(figsize=(40/2.54, 40/2.54))
446 frame1 = fig1.add_subplot(1,1,1,xscale="log",yscale="log")
447 frame1.set_title("{}".format(names[n]),fontsize=30)
448 frame1.set_xlim([0.25,30])
449 frame1.errorbar(v,S,yerr=Serr,ecolor="k",marker="*",mfc="y",ms=10,ls="",label="Recorded
450 data")
450 if q == 0 or q == 3 or q == 5:
451     print("Power law has no peak: no peak has been plotted.")
452 else:
453     frame1.errorbar(vmax,Smax,xerr=vmaxerr,yerr=Smaxerr,ecolor="k",marker="o",mfc="b",
454 ms=15,ls="",alpha=0.5,label="Peak")
454 frame1.plot(vfit,fitfun,"b",label="Fitted function (least squares)")
455 if q == 2:
456     frame1.plot(vfit,Sfitsyn(vfit,p[0],p[1],p[2],p[3]),"r",label="Fitted function, power
457 lawless")
457 frame1.fill_between(vfit,fitmax,fitmin,facecolor="blue",alpha=0.15)
458 frame1.set_xlabel('Frequency v (GHz)',fontsize=25)
459 frame1.set_ylabel('Flux density S (mJy)',fontsize=25)
460 frame1.tick_params(labelsize=25)
461
462 ###Plotting 2d multivariate pdf results (Episode VI).
463 fig2 = figure(figsize=(40/2.54, 40/2.54))
464 if q == 0:
465     a1,a2,pospos = mpdfax(p[0],p[1],pcov[0,0],pcov[1,1])
466     frame2 = fig2.add_subplot(1,1,1)
467     frame2.contourf(a1,a2,mpdf(p[0],p[1],pcov[0,0],pcov[0,1],pcov[1,0],pcov[1,1],pospos)
468 )
469     frame2.set_xlabel('coefficient ${}$'.format(coeflist[0+os]))
470     frame2.set_ylabel('coefficient ${}$'.format(coeflist[1+os]))
471 elif q == 1 or q == 5:
472     k = 0
473     for i in range(ndim):
474         for j in range(ndim):
475             if j > i:
476                 a1,a2,pospos = mpdfax(p[i],p[j],pcov[i,i],pcov[j,j])
477                 frame2 = fig2.add_subplot(2,3,k+1)
478                 k = k + 1
479                 frame2.contourf(a1,a2,mpdf(p[i],p[j],pcov[i,i],pcov[i,j],pcov[j,i],pcov[
480 j,j],pospos))
481                 frame2.set_xlabel('coefficient ${}$'.format(coeflist[i+os]))
482                 frame2.set_ylabel('coefficient ${}$'.format(coeflist[j+os]))
483 elif q == 2:
484     print("Multivariate pdf's could not be plotted because of the limitations of
485 detestable frequentist statistics.")
486 elif q == 3:
487     a1,a2,pospos = mpdfax(p[0],p[1],pcov[0,0],pcov[1,1])
488     frame2 = fig2.add_subplot(1,1,1)
489     frame2.contourf(a1,a2,mpdf(p[0],p[1],pcov[0,0],pcov[0,1],pcov[1,0],pcov[1,1],pospos)
490 )

```

```

487     frame2.set_xlabel('coefficient ${}$'.format(coeflist[0+os]))
488     frame2.set_ylabel('coefficient ${}$'.format(coeflist[1+os]))
489 elif q == 4:
490     k = 0
491     for i in range(ndim):
492         for j in range(ndim):
493             if j > i:
494                 a1,a2,pospos = mpdfax(p[i],p[j],pcov[i,i],pcov[j,j])
495                 frame2 = fig2.add_subplot(3,4,k+1)
496                 k = k + 1
497                 frame2.contourf(a1,a2,mpdf(p[i],p[j],pcov[i,i],pcov[i,j],pcov[j,i],pcov[
j,j],pospos))
498                 frame2.set_xlabel('coefficient ${}$'.format(coeflist[i+os]))
499                 frame2.set_ylabel('coefficient ${}$'.format(coeflist[j+os]))
500
501 show()
502
503 while True:
504     print("\033[1m","Do you wish to save the data and figures?","\033[0m")
505     answer = input(" :")
506     if answer.lower().startswith("y"):
507         print("Roger")
508         break
509     elif answer.lower().startswith("n"):
510         print("OKDOEP")
511         exit()
512
513 ###Saving figures to folder.
514 fig1.savefig("/Users/users/tjoa/Downloads/Thesis/Pics/Part1/Synchfits/Synfit{0}-{1}.png"
.format(n,q))
515 if q == 2:
516     print("No multivariate pdf's were made and thus none were saved to directory.")
517 else:
518     fig2.savefig("/Users/users/tjoa/Downloads/Thesis/Pics/Part1/Contourplots/Conplt{0}-
{1}.png".format(n,q))
519
520 ###Writing values to file.
521 if q == 0:
522     with open("ThesisTable_powerlaw0.dat", "a") as myfile:
523         myfile.write("{0} {1} {2} {3} {4} {5} {6}" .format(n,names[n],p[0],pcov[0,0],p
[1],pcov[1,1],"regpls"))
524         myfile.write("\n")
525 elif q == 1:
526     with open("ThesisTable_synchrotron0.dat", "a") as myfile:
527         myfile.write("{0} {1} {2} {3} {4} {5} {6} {7} {8} {9} {10} {11} {12} {13}" .
format(n,names[n],p[0],pcov[0,0],p[1],pcov[1,1],p[2],pcov[2,2],p[3],pcov[3,3],vmax,
vmaxerr,Smax,Smaxerr,))
528         myfile.write("\n")
529 elif q == 2:
530     with open("ThesisTable_synch-powlaw0.dat", "a") as myfile:
531         myfile.write("{0} {1} {2} {3} {4} {5} {6} {7} {8} {9} {10} {11} {12} {13} {14}
{15} {16} {17} {18}" .format(n,names[n],p[0],pcov[0,0],p[1],pcov[1,1],p[2],pcov[2,2],
p[3],pcov[3,3],p[4],pcov[4,4],p[5],pcov[5,5],vmax,vmaxerr,Smax,Smaxerr,"MCMC"))
532         myfile.write("\n")
533 elif q == 3:
534     with open("ThesisTable_powerlaw0.dat", "a") as myfile:
535         myfile.write("{0} {1} {2} {3} {4} {5} {6}" .format(n,names[n],p[0],pcov[0,0],p
[1],pcov[1,1],"agepls"))
536         myfile.write("\n")
537 elif q == 4:
538     with open("ThesisTable_synch-powlaw0.dat", "a") as myfile:
539         myfile.write("{0} {1} {2} {3} {4} {5} {6} {7} {8} {9} {10} {11} {12} {13} {14}
{15} {16} {17}" .format(n,names[n],p[0],pcov[0,0],p[1],pcov[1,1],p[2],pcov[2,2],p[3],
pcov[3,3],p[4],pcov[4,4],M,0,vmax,vmaxerr,Smax,Smaxerr))
540         myfile.write("\n")
541 elif q == 5:
542     with open("ThesisTable_double-powlaw0.dat", "a") as myfile:
543         myfile.write("{0} {1} {2} {3} {4} {5} {6} {7} {8} {9}" .format(n,names[n],p[0],
pcov[0,0],p[1],pcov[1,1],p[2],pcov[2,2],p[3],pcov[3,3]))
544         myfile.write("\n")

```

## B.2 Markov-chain Monte Carlo fitting algorithm

The following script is used to fit SSA GPS+power law spectra with fixed index  $m$  using an MCMC algorithm.

Scripts/Thesis\_Sv-MCMCfit\_synpl14-3.py

```
1 #!/usr/bin/env python
2 from __future__ import division
3 import numpy as np
4 import scipy.optimize as spo
5 import emcee as mc
6 import corner as co
7 from matplotlib.pyplot import figure, show
8 #####THIS SCRIPT CAN RUN MCMC ON SYNCHROTRON PLUS POWER LAW SPECTRA ONLY, FOR FIXED m.
9
10 #####Episode I: Reading Data (or: The Phantom Data).
11 #####
12 print(">>> Episode I: The Phantom Data")
13
14 #####Reading relevant data from (mono-axial) datafile. If uncertainties were unknown, they
15 have been set to 1.0 to prevent code collapse due to divide by zero errors.
16 fulllist = np.genfromtxt("Janskylist.dat", dtype=None, usecols=[0], comments="#")
17 errlist = np.genfromtxt("Janskylist.dat", dtype=None, usecols=[1], comments="#")
18
19 #####Splitting the superarray into useful, separate arrays with relevant data.
20 names, S003, S014, S048, S084, S149, S224 = np.split(fulllist, 7)
21 junkerr, errS003, errS014, errS048, errS084, errS149, errS224 = np.split(errlist, 7)
22 names = names.astype(str).flatten()
23 S003 = np.asarray(S003.astype(float).flatten())
24 S014 = np.asarray(S014.astype(float).flatten())
25 S048 = np.asarray(S048.astype(float).flatten())
26 S084 = np.asarray(S084.astype(float).flatten())
27 S149 = np.asarray(S149.astype(float).flatten())
28 S224 = np.asarray(S224.astype(float).flatten())
29
30 #####Printing all data in a clear-cut table.
31 print("\033[1m", "index * name * 0.325 * 1.4 * 4.85 * 8.46 * 14.94 * 22.46", "\033[0m",
32 <- Frequency bands")
33 for i in range(len(names)):
34     print(i, ":", "\033[1m", names[i], "\033[0m", "*", S003[i], "*", S014[i], "*", S048[i], "*",
35           S084[i], "*", S149[i], "*", S224[i])
36 n = int(input(" Pick GPS source: "))
37
38 #####Creating the to be used data arrays.
39 v000 = [0.325, 1.4, 4.85, 8.46, 14.94, 22.46]
40 S000 = [S003[n], S014[n], S048[n], S084[n], S149[n], S224[n]]
41 Serr000 = [errS003[n], errS014[n], errS048[n], errS084[n], errS149[n], errS224[n]]
42
43 v = np.asarray(v000)
44 S = np.asarray(S000)
45 Serr = np.asarray(Serr000)
46
47 print("Source designation:", names[n])
48 print("Frequency bands:", v)
49 print("Flux densities:", S)
50 print("Flux uncertainties:", Serr)
51
52 #####Plotting the recorded data points for visualisation purposes.
53 fig = figure(figsize=(10,10))
54 frame = fig.add_subplot(1,1,1, xscale="log", yscale="log")
55 frame.errorbar(v, S, yerr=Serr, ecolor="k", marker="*", mfc="y", ms=10, ls="", label="Data from
56 {}" .format(names[n]))
57 frame.set_xlabel('Frequency v (GHz)')
58 frame.set_ylabel('Flux density S (mJy)')
59 frame.legend()
60 show()
61
62 #####While-loop to allow user-friendly cancelling of code.
63 while True:
```

```

59 print("\033[1m","Do you want to continue?","\033[0m")
60 answer = input(" :")
61 if answer.lower().startswith("y"):
62     print("Roger: synch fit and power law online")
63     break
64 elif answer.lower().startswith("n"):
65     print("OKDOEF")
66     exit()
67
68 ###Setting model type to synchrotron plus power law.
69 q = 2 ; ndim = 5 ; os=0
70 modelname = ["power law","synchrotron","synchrotron/power law"]
71 print(" Number of free parameters:",ndim)
72 M = (np.log10(0.5*S[1])-np.log10(S[0]))/(np.log10(v[1])-np.log10(v[0]))
73
74
75 #####Episode II: Creating probability functions (or: Attack of the Functions).
76 #####
77 print(">>> Episode II: Attack of the Functions")
78
79 def Sfitsyn(v,S01,v01,k,l):
80     f1 = S01*(v/v01)**k          ###First power law
81     f2 = (1-np.e**(-(v/v01)**(1-k)))/(1-np.e**(-1))    ###Second power law (exponential
82     fall off)
83     return f1*f2
84
85 def Sfitplsyn(v,S01,v01,k,l,A):
86     f1 = S01*(v/v01)**k          ###First power law
87     f2 = (1-np.e**(-(v/v01)**(1-k)))/(1-np.e**(-1))    ###Second power law (exponential
88     fall off)
89     f3 = A*(v)**M                ###Third power law (only when two turnovers
90     are present)
91     return f1*f2 + f3
92
93 def errSfitplsyn(v,S01,v01,k,l,A,cov):    ###Finding uncertainties function in synch fit
94     plus power law.
95     C = S01/(np.e-1)
96     C2 = S01/(1-1/np.e)
97     E = np.e**(-(v/v01)**(1-k))
98     E2 = np.e**((1-((v/v01)**(1-k))))
99     dsdS01 = Sfitsyn(v,S01,v01,k,l)/S01
100     dsdv01 = (-((C2*(1-E)))*k*v**((v/v01)**(k-1)))/(v01**2) - (C2*(1-k)*v**E**((v/v01)**(1
101     -1)))/(v01**2)
102     dsdk = C2*(1-E)*((v/v01)**k)*np.log(v/v01) - C2*((v/v01)**(1))*np.log(v/v01)*E
103     dsdl = C*((v/v01)**1)*np.log(v/v01)*E2
104     dsdA = v**M
105     dsdm = A*(v**M)*np.log(v)
106     dsdx = np.asarray([dsdS01,dsdv01,dsdk,dsdl,dsdA,dsdm])
107     Svar = 0
108     for i in range(ndim):
109         for j in range(ndim):
110             Svar = Svar + dsdx[i]*dsdx[j]*cov[i,j]
111     return np.sqrt(Svar)
112
113 def derSfitplsyn(v,S01,v01,k,l,A):    ###Derivative of power law/synchrotron
114     combined fit
115     E = np.e**(-(v/v01)**(1-k))
116     C = S01/(v01*(1-1/np.e))
117     f1 = C*k*((v/v01)**(k-1))*(1-E)
118     f2 = C*(1-k)*((v/v01)**(1-1))*E
119     f3 = A*M*v**(M-1)
120     return f1 + f2 + f3
121
122 ###Array spanning the fitting space.
123 vfit = np.linspace(0.3,23,1000)
124
125 ###List that names the coefficients for the synchfit.
126 coeflist = ["S_{01}","v_{01}","k","l","A","m"]

```

```

122 #####Episode III: Computing estimates for free parameters. (or: Revenge of the Estimates
    ). #####
123 print(">>> Episode III: Revenge of the Estimates")
124
125 ###Creating some new, useful arrays and indices.
126 S2 = [S[1],S[2],S[3],S[4],S[5]]
127 v2 = [v[1],v[2],v[3],v[4],v[5]]
128 Smi = S000.index(max(S000))
129 Smi2 = S2.index(max(S2))
130
131 ###Finding estimates for free parameters.
132 S0guess = S[4]
133 v0guess = v[4]
134 if S2[1] > S2[0]:
135     kguess = (np.log10(S2[1])-np.log10(0.5*S2[0]))/(np.log10(v2[1])-np.log10(v2[0]))
136 else:
137     if S2[2] > S2[1]:
138         kguess = (np.log10(S2[2])-np.log10(0.5*S2[1]))/(np.log10(v2[2])-np.log10(v2[1]))
139     else:
140         kguess = 2.0
141 if S2[4] < S2[3]:
142     lguess = (np.log10(S2[4])-np.log10(S2[3]))/(np.log10(v2[4])-np.log10(v2[3]))
143 else:
144     if S2[4] < S2[2]:
145         lguess = (np.log10(S2[4])-np.log10(S2[2]))/(np.log10(v2[4])-np.log10(v2[2]))
146     else:
147         lguess = -1.0
148 Aguess = S[0]/(v[0]**M)
149
150 estimates = [S0guess, v0guess, kguess, lguess, Aguess, M]
151 print("Parameter estimates:")
152 print(np.asarray(estimates))
153
154 pg, pgcov = spo.curve_fit(Sfitplsyn, v, S, [S0guess, v0guess, kguess, lguess, Aguess], sigma=
    Serr, bounds=([0,0.3,0,-2.5,0],[500,23,2.5,0,100]),**{'max_nfev':100000})
155 for i in range(ndim-1):
156     print(pg[i], pgcov[i, i])
157 print(M)
158
159 Sstd = 2.0 ; Slow = 0.0 ; Supp = 500.0 #####Constraints on S0
160 vstd = 1.0 ; vlow = v[1] ; vupp = 23.0 #####Constraints on v0
161 kstd = 0.5 ; klow = 0 ; kupp = 2.5 #####Constraints on k
162 lstd = 0.5 ; llow = -2.5 ; lupp = 0.0 #####Constraints on l
163 Astd = 2.0 ; Alow = 0.0 ; Aupp = 100.0 #####Constraints on A
164
165 def normdist(x, mean, std): #####Normal (Gaussian) distribution
166     expterm = ((x-mean)**2)/(2*std**2)
167     f = (1/(std*np.sqrt(2*np.pi)))*np.e**-expterm
168     return f
169
170 def lnprior(theta): #####ln of prior
171     S01, v01, k, l, A = theta
172     if Slow < S01 < Supp and vlow < v01 < vupp and klow < k < kupp and llow < l < lupp
    and Alow < A < Aupp:
173         return np.log(normdist(k, pg[2], kstd)*normdist(l, pg[3], lstd))
174     return -np.inf
175
176 def lnlike(theta, v, S, Serr): #####ln of likelihood
177     S01, v01, k, l, A = theta
178     f1 = S01*(v/v01)**k #####First power law
179     f2 = (1-np.e**(-(v/v01)**(1-k)))/(1-np.e**(-1)) #####Second power law (exponential
    fall off)
180     f3 = A*(v)**M #####Third power law (only when two turnovers are
    present)
181     model = f1*f2+f3
182     inv_sigma2 = 1.0/(Serr**2)
183     return -0.5*(np.sum((S-model)**2*inv_sigma2 + np.log(2*np.pi/inv_sigma2)))
184
185 def lnprob(theta, v, S, Serr): #####ln of posterior
186     lp = lnprior(theta)

```

```

187     if lp == -np.inf:
188         return -np.inf
189     return lp + lnlike(theta, v, S, Serr)
190
191
192 #####Episode IV: Performing MCMC. (or: A New Method).
193 #####
194 print(">>> Episode IV: A New Method")
195
196 ###Setting dimensions, walkers and mainframe.
197 nwalkers = 300
198 p0 = [pg + 1e-2*np.random.randn(ndim) for i in range(nwalkers)]
199 sampler = mc.EnsembleSampler(nwalkers, ndim, lnprob, args=(v, S, Serr))
200
201 ###Burn-in & reset.
202 print(" Performing burn-in sequence...")
203 pos, prob, state = sampler.run_mcmc(p0, 200)
204 sampler.reset()
205
206 #If lnprob0 and newlnprob BOTH return -inf, then lnpdiff is nan. If lnprob0 ALONE
207 #returns -inf, lnpdiff is inf. If newlnprob ALONE returns -inf, lnpdiff is -inf.
208
209 ###Running real MCMC!
210 print(" Initiating main MCMC algorithm...")
211 nsteps = 1000
212 sampler.run_mcmc(pos, nsteps, rstate0=state)
213
214 ###Generating the cornerplots
215 print(" Generating cornerplots...")
216 burnin = 50
217 samples = sampler.chain[:, burnin:, :].reshape((-1, ndim))
218
219 #####Episode V: Finding parameter value & covariance matrix (or: The Parameters Strike
220 #Back). #####
221 print(">>> Episode V: The Parameters Strike Back")
222
223 resolution = 1000
224
225 def MCs(a, i): #####Sampler function.
226     return a.flatchain[:, i]
227
228 def histo(x): #####Histogram maker.
229     hist = np.histogram(x, resolution)
230     counts = hist[0] ; values = hist[1]
231     values = np.delete(values, resolution)
232     a = np.asarray([])
233     for i in range(resolution):
234         appa = np.full(counts[i], values[i])
235         a = np.append(a, appa)
236     return a, counts, values
237
238 ###Creating total probability array from six dimensions, extracting maximum probability
239 #index.
240 lnprobarr = sampler.flatlnprobability
241 mindex = np.argmax(lnprobarr)
242
243 ###Filling the parameter array and the covariance matrix.
244 phist = []
245 p = np.zeros(ndim)
246 print("\033[1m+" >>> Using MCMC on {} model:".format(modelname[q]), "\033[0m")
247 for i in range(ndim):
248     hishis = histo(MCs(sampler, i))
249     phist.append(hishis[0])
250     p[i] = MCs(sampler, i)[mindex]
251
252 phist = np.asarray(phist)
253 pcov = np.cov(phist)
254
255 for i in range(ndim):

```

```

253     print(" Coefficient",coeflist [i+os],"equals",p[i])
254     print("   Uncertainty of coefficient",coeflist [i+os],"equals",np.sqrt(pcov[i,i]))
255 print(" Coefficient m equals:",M)
256
257
258 #####Episode VI: Finding the maximum of the fit (or: Return of the Maxima).
259 #####
260 print(">>> Episode VI: Return of the Maxima")
261
262 ###Array spanning the bisection space, hence excluding the minimum to get the turnover
263 as return from scipy.bisect.
264 if S[2] > S[1]:
265     inc = (v[2]+v[1])/2
266 else:
267     if S[3] > S[2]:
268         inc = (v[3]+v[2])/2
269     else:
270         inc = v[3]
271
272 vfitbis = np.linspace(inc,23,1000)
273
274 def bisector(v):
275     #####Function of one variable used by the bisection
276     algorithm.
277     return derSfitplsyn(v,*p)
278
279
280 #####If loop that bisects the function if it has one or more peaks.
281 for i in range(len(vfitbis)-1):
282     if bisector(vfitbis [i+1]) < 0 and bisector(vfitbis [i]) > 0:
283         ma = vfitbis [i+1]
284         mi = vfitbis [i]
285 vmax = spo.bisect(bisector,mi,ma)
286 vmaxerr = (23-inc)/1000 #Range and stepsize of fitting space.
287 Smax = Sfitplsyn(vmax,*p)
288 Smaxerr = errSfitplsyn(vmax,*p,pcov)
289
290 #####Printing peak coordinates.
291 print("\033[1m", "   >>> Using scipy.optimize.bisect on {} model gives peak coordinates:
292       ".format(modelname[q]), "\033[0m")
293 print(" Fit frequency of peak is:",vmax,"GHz")
294 print("   Uncertainty of frequency peak is:",vmaxerr,"GHz")
295 print(" Fit flux density of peak is:",Smax,"mJy")
296 print("   Uncertainty of flux peak is:",Smaxerr,"mJy")
297
298
299 #####Episode VII: Plotting all results (or: The Pyplot Awakens).
300 #####
301 print(">>> Episode VII: The Pyplot Awakens")
302
303 fitfun = Sfitplsyn(vfit,*p)
304 fitfun2 = Sfitplsyn(vfit,*pg)
305 fitmax = fitfun + errSfitplsyn(vfit,*p,pcov)
306 fitmin = fitfun - errSfitplsyn(vfit,*p,pcov)
307
308
309 #####Plotting data, least squares fits and peaks (Episodes I, IV & V).
310 fig1 = figure(figsize=(40/2.54, 40/2.54))
311 frame1 = fig1.add_subplot(1,1,1,xscale="log",yscale="log")
312 frame1.set_title("{} {}".format(names[n]), fontsize=30)
313 frame1.set_xlim([0.25,30])
314 frame1.errorbar(v,S,yerr=Serr,ecolor="k",marker="*",mfc="y",ms=10,ls="",label="Recorded
315 data")
316 frame1.errorbar(vmax,Smax,xerr=vmaxerr,yerr=Smaxerr,ecolor="k",marker="o",mfc="b",ms
317 =15,ls="",alpha=0.5,label="Peak")
318 frame1.plot(vfit,fitfun,"b",label="Fitted function (MCMC-max)")
319 frame1.plot(vfit,fitfun2,"r",label="Fitted function (least squares)")
320 frame1.fill_between(vfit,fitmax,fitmin,facecolor="blue",alpha=0.15)
321 frame1.set_xlabel('Frequency v (GHz)', fontsize=25)
322 frame1.set_ylabel('Flux density S (mJy)', fontsize=25)
323 frame1.tick_params(labelsize=25)
324
325
326 #####Plotting cornerplots

```

```

316 fig2 = co.corner(samples, labels=["$$_{01}$", "$v_{01}$", "$k$", "$l$", "$A$", "$m$"])
317
318 ###Plotting MCMC results (histograms for each coefficient) (Episode V).
319 fig3 = figure(figsize=(40/2.54, 40/2.54))
320 for i in range(ndim):
321     frame3 = fig3.add_subplot(3,2,i+1)
322     frame3.hist(MCs(sampler, i), normed=True, bins=200, alpha=0.8)
323     frame3.axvline(x=p[i], linewidth=2, color='r')
324     frame3.axvline(x=pg[i], linewidth=1, color='g')
325     frame3.set_xlabel("coefficient {}".format(coeflist[i+os]))
326     frame3.set_ylabel("walker counts")
327
328 show()
329
330 while True:
331     print("\033[1m", "Do you wish to save the data and figures?", "\033[0m")
332     answer = input(" :")
333     if answer.lower().startswith("y"):
334         print("Roger")
335         break
336     elif answer.lower().startswith("n"):
337         print("OKDOE!")
338         exit()
339
340 ###Saving figures to folder.
341 fig1.savefig("/Users/users/tjoa/Downloads/Thesis/Pics/Part1/Synchfits/Synfit{0}-{1}-{2}.
    png".format(names[n], n, q))
342 fig2.savefig("/Users/users/tjoa/Downloads/Thesis/Pics/Part1/Contourplots/Conplt{0}-{1}-
    {2}.png".format(names[n], n, q))
343 fig3.savefig("/Users/users/tjoa/Downloads/Thesis/Pics/Part1/MCMCHistograms/Histogram
    {0}-{1}-{2}.png".format(names[n], n, q))
344
345
346 ###Writing values to file.
347 with open("ThesisTable.synch-powlaw4.dat", "a") as myfile:
348     myfile.write("{0} {1} {2} {3} {4} {5} {6} {7} {8} {9} {10} {11} {12} {13} {14} {15}
    {16} {17}".format(n, names[n], p[0], pcov[0,0], p[1], pcov[1,1], p[2], pcov[2,2], p[3], pcov
    [3,3], p[4], pcov[4,4], M, "N/A", vmax, vmaxerr, Smax, Smaxerr))
349     myfile.write("\n")

```

### B.3 K-z relation fitting algorithm

The following script is used to fit the K-z relation and obtain redshifts from our sample.

Scripts/Thesis\_Kz-relation\_v13-1.py

```

1 #!/usr/bin/env python
2 from __future__ import division
3 import numpy as np
4 import scipy.stats as sps
5 import numpy.random as npr
6 from matplotlib.pyplot import figure, show
7 ###THIS SCRIPT COMPUTES STD IN LOG10(z) AND z FROM K-BAND MAGNITUDES.
8
9
10 #####Episode I: Reading Data (or: The Phantom Data).
11 #####
12 print(">>> Episode I: The Phantom Data")
13
14 ###Reading relevant data from datafile.
15 datanames = np.genfromtxt("fulltable.dat", dtype=str, usecols=[0], comments="#")
16 zlist = np.genfromtxt("fulltable.dat", dtype=float, usecols=[1], comments="#")
17 zstdlowlist = np.genfromtxt("fulltable.dat", dtype=float, usecols=[2], comments="#")
18 zstdupplist = np.genfromtxt("fulltable.dat", dtype=float, usecols=[3], comments="#")
19 Klist = np.genfromtxt("fulltable.dat", dtype=float, usecols=[4], comments="#")
20 Kstdlist = np.genfromtxt("fulltable.dat", dtype=float, usecols=[5], comments="#")
21
22 ###Printing all data in a clear-cut table.

```

```

22 print("\033[1m", "index * source * z *zstdlow *zstdup * K-mag * K-std"\033[0m"
    )
23 for i in range(len(datanames)):
24     print(i, "      :", "\033[1m", datanames[i], "\033[0m", "{0} {1} {2} {3} {4}".
        format(zlist[i], zstdlowlist[i], zstduplist[i], Klist[i], Kstdlist[i]))
25
26 #####Episode II: Defining the functions and general constants (or: Attack of the
    Functions). #####
27 print(">>> Episode II: Attack of the Functions")
28
29 ###Function block.
30 def Klogzrel(logz):          ###K,log10(z)-relation function
31     return 17.37 + 4.53*logz + 0.31*logz**2
32
33 def Kzrel(z):                ###K,z-relation function
34     return 17.37 + 4.53*np.log10(z) + 0.31*(np.log10(z))**2
35
36 ###Specifying general constants and limits of the fit.
37 stdK1 = 0.593
38 logzmin = -4 ; logzmax = np.log10(10)
39 ranlen = 5000000
40
41
42
43 #####Episode III: Generating fit line and random data (or: Revenge of the Randomness).
    #####
44 print(">>> Episode III: Revenge of the Randomness")
45
46 ###Generating plain , idealized K-log10(z) relation.
47 logzarr = np.linspace(logzmin, logzmax, ranlen)
48 Karr = Klogzrel(logzarr)
49
50 ###Generating the random points around idealized K-log10(z) relation.
51 logzran = (logzmax-logzmin)*npr.random(ranlen)+logzmin
52 Kran = np.zeros(ranlen)
53 for i in range(ranlen):
54     Kran[i] = npr.normal(Klogzrel(logzran[i]), stdK1)
55
56
57 #####Episode IV: Finding redshift uncertainties in data (or: A New Data).
    #####
58 print(">>> Episode IV: A New Data")
59
60 def uberfunctionarr(DATA, DATAerr): ###Function that computes distribution around an
    array of Ksamples
61     ###Specifying arrays and values to be used in log10(z)/z uncertainty computation
    loop as empty zero arrays.
62     Ksample = DATA          ###Data sample of K points
63     Kdataerr = DATAerr      ###K data uncertainties
64     dic = {}                 ###Empty dictionary for log10(z) values, to be
        filled in loop
65     dic2 = {}                ###Empty dictionary for z values, to be filled
        in loop
66     logzsample = np.zeros(len(Ksample))    ###Sample of log10(z) points
67     zsample = np.zeros(len(Ksample))       ###Corresponding sample of z points
68     stdK = stdK1*np.ones(len(Ksample))     ###Array of std_K values of sample
        length
69     stdlogz = np.zeros(len(Ksample))       ###Empty std_log10(z) array
70     stdzlow = np.zeros(len(Ksample))       ###Empty std_z (lower bound) array
71     stdzupp = np.zeros(len(Ksample))       ###Empty std_z (upper bound) array
72     step = 0.05                        ###Step size for the loop
73
74     ###Loop that computes the uncertainty in log10(z) and z at some z using an algorithm
        that takes samples from a small interval, then creates the corresponding
        distribution and returns the std in log10(z) and z.
75     for i in range(len(Ksample)):
76         Kupp = Ksample[i]+step ; Klow = Ksample[i]-step
            ###Upper and lower interval bounds
77         dic["logzdist-{}".format(i)] = np.asarray([])
            ###Creating new log10(z) dictionary entry

```

```

78     dic2["zdist-{}".format(i)] = np.asarray([])
79     ###Creating new z dictionary entry
80     for j in range(ranlen):
81         if Kran[j] > Klow and Kran[j] < Kupp:
82             ###Acceptance condition
83             dic["logzdist-{}".format(i)] = np.append(dic["logzdist-{}".format(i)],
84             logzran[j])             ###Appending accepted values to corresponding log10(z) dictionary
85             entry
86             dic2["zdist-{}".format(i)] = np.append(dic2["zdist-{}".format(i)],10**
87             logzran[j])             ###Appending accepted values to corresponding z dictionary entry
88             data = dic["logzdist-{}".format(i)]
89             ###Renaming dictionary entry for easy handling
90             data2 = dic2["zdist-{}".format(i)]
91             ###Renaming dictionary entry for easy handling
92
93             ###Fitting a normal distribution to dictionary data.
94             x = np.linspace(data.min(), data.max(), 100)             ###Creating regular
95             possible value space
96             param = sps.norm.fit(data)             ###Fitting normal parameters
97             to the data with the mean at the corresponding log10(z) value.
98             pdf_fitted = sps.norm.pdf(x, *param)             ###Fitting a normal
99             distribution with said parameters to the interval
100             prob = pdf_fitted/pdf_fitted.sum()             ###Normalizing pdf
101
102             ###Fitting a log-normal distribution to dictionary 2 data.
103             x2 = np.linspace(data2.min(), data2.max(), 100)             ###Creating
104             regular possible value space
105             param2 = sps.lognorm.fit(data2)             ###Fitting log-normal
106             parameters to the data with the mean at the corresponding z value.
107             pdf_fitted2 = sps.lognorm.pdf(x2, *param2)             ###Fitting a log-
108             normal distribution with said parameters to the interval
109             prob2 = pdf_fitted2/pdf_fitted2.sum()             ###Normalizing pdf
110
111             ###Computing required statistics from dictionary data.
112             logzmu = x.dot(prob)             ###Computing mean in log10(z)
113             )
114             logzsample[i] = logzmu             ###Entering into ex-
115             loop logz array
116             zmu = 10**logzmu             ###Computing mean in z
117             zsample[i] = zmu             ###Entering into ex-loop z
118             array
119             stdlogz1 = np.sqrt(np.power(x,2).dot(prob) - logzmu**2)             ###Computing std
120             in log10(z) (sqrt(data_logz^2 - mu_logz^2))
121             stdlogz[i] = stdlogz1             ###Entering into ex-loop
122             stdlogz array
123             stdzlow1 = 10**(logzmu-stdlogz1)-zmu             ###Computing std lower
124             bound in z 10^(mu_logz-std_logz)-mu_z
125             stdzlow[i] = -1*stdzlow1             ###Entering into ex-loop
126             stdzlow array
127             stdzupp1 = 10**(logzmu+stdlogz1)-zmu             ###Computing std upper
128             bound in z 10^(mu_logz+std_logz)-mu_z
129             stdzupp[i] = stdzupp1             ###Entering into ex-loop
130             stdzipp array
131             return (Ksample, Kdataerr, logzsample, stdlogz, zsample, stdzlow, stdzupp, dic, dic2)
132
133 def uberfunctionflt (DATA,DATAerr): ###Function that computes distribution around a
134     single Ksample
135     ###Specifying arrays and values to be used in log10(z)/z uncertainty computation
136     loop as empty zero arrays.
137     Ksample = DATA             ###Data sample of K points
138     Kdataerr = DATAerr             ###K data uncertainties
139     stdK = stdK1             ###Array of std_K values of sample length
140     logzdist = np.asarray([])             ###Set of logz-values
141     zdist = np.asarray([])             ###Set of z-values
142     step = 0.05             ###Step size
143
144     ###Computing the uncertainty in log10(z) and z at some z using an algorithm that
145     takes samples from a small interval, then creating the corresponding distribution
146     and returning the std in log10(z) and z.
147     Kupp = Ksample+step ; Klow = Ksample-step             ###Upper and lower

```

```

122 interval bounds
123 for j in range(ranlen):
124     if Kran[j] > Klow and Kran[j] < Kupp:                                     ###Acceptance
125     condition
126         logzdist = np.append(logzdist, logzran[j])                             ###Appending accepted
127         values to corresponding log10(z) dictionary entry
128         zdist = np.append(zdist, 10**logzran[j])                               ###Appending accepted
129         values to corresponding z dictionary entry
130         data = logzdist                                                         ###Renaming dictionary entry
131         for easy handling
132         data2 = zdist                                                            ###Renaming dictionary entry for
133         easy handling
134
135     ###Fitting a normal distribution to data.
136     x = np.linspace(data.min(), data.max(), 100)                               ###Creating regular
137     possible value space
138     param = sps.norm.fit(data)                                                  ###Fitting normal
139     parameters to the data with the mean at the corresponding log10(z) value.
140     pdf_fitted = sps.norm.pdf(x, *param)                                       ###Fitting a normal
141     distribution with said parameters to the interval
142     prob = pdf_fitted/pdf_fitted.sum()                                         ###Normalizing pdf
143
144     ###Fitting a log-normal distribution to data2.
145     x2 = np.linspace(data2.min(), data2.max(), 100)                           ###Creating regular
146     possible value space
147     param2 = sps.lognorm.fit(data2)                                            ###Fitting log-normal
148     parameters to the data with the mean at the corresponding z value.
149     pdf_fitted2 = sps.lognorm.pdf(x2, *param2)                                ###Fitting a log-
150     normal distribution with said parameters to the interval
151     prob2 = pdf_fitted2/pdf_fitted2.sum()                                     ###Normalizing pdf
152
153     ###Computing required statistics from data.
154     logzmu = x.dot(prob)                                                         ###Computing mean in log10(z)
155     )
156     zmu = 10**logzmu                                                            ###Computing mean in z
157     stdlogz1 = np.sqrt(np.power(x,2).dot(prob) - logzmu**2)                   ###Computing std in
158     log10(z) (sqrt(data_logz^2 - mu_logz^2))
159     stdzlow1 = 10**(logzmu-stdlogz1)-zmu                                       ###Computing std lower bound
160     in z 10^(mu_logz-std_logz)-mu_z
161     stdzuppl = 10**(logzmu+stdlogz1)-zmu                                       ###Computing std upper bound
162     in z 10^(mu_logz+std_logz)-mu_z
163     return (Ksample, Kdataerr, logzmu, stdlogz1, zmu, stdzlow1, stdzuppl)
164
165 newKlist = []
166 newKerrlist = []
167 newdatanames = []
168 for i in range(len(Klist)):
169     if Klist[i] == 0.0:
170         print("Entry", i, "has no known K-band magnitude.")
171     else:
172         UBER = uberfunctionflt(Klist[i], Kstdlist[i])
173         zlist[i] = UBER[4]
174         zstdlowlist[i] = np.abs(UBER[5])
175         zstduppllist[i] = UBER[6]
176         newKlist.append(Klist[i])
177         newKerrlist.append(Kstdlist[i])
178         newdatanames.append(datanames[i])
179 newKarr = np.asarray(newKlist)
180 newKerrarr = np.asarray(newKerrlist)
181 fdata5 = uberfunctionarr(newKarr, newKerrarr)
182
183 ###Printing all data in a clear-cut table.
184 print("\033[1m", "index * source * z * zstdlow * zstdup * K-mag * K-std" "\033[0m"
185 )
186 for i in range(len(datanames)):
187     print(i, " ", " ", "\033[1m", datanames[i], "\033[0m", "{0} {1} {2} {3} {4}".
188     format(np.round(zlist[i], 4), np.round(zstdlowlist[i], 4), np.round(zstduppllist[i], 4),
189     Klist[i], Kstdlist[i]))
190
191
192

```

```

173 #####Episode V: Plotting results (or: The Pyplot Strikes Back).
174 #####
175 print(">>> Episode V: The Pyplot Strikes Back")
176
177 ###Plot no.1: K-band magnitude versus log10(z) via idealized K,log10(z)-relation and
178 simulated data around it.
179 fig1 = figure(figsize=(40/2.54, 40/2.54))
180 frame1 = fig1.add_subplot(1,1,1)
181 frame1.set_title("K-band magnitude vs log10(z)", fontsize=15)
182 frame1.plot(logzran ,Kran, 'g.', alpha=0.1, label="Simulated random distribution")
183 frame1.plot(logzarr ,Karr, linewidth=5, alpha=0.5, label="Idealized K,log10(z)-relation")
184 frame1.errorbar(fdata5[2], fdata5[0], xerr=fdata5[3], yerr=fdata5[1], ecolor="k", elinewidth
185 =3, marker="*", mfc="y", ms=10, ls="") , label="Sample K, z")
186 frame1.set_xlabel('Redshift log10(z)')
187 frame1.set_ylabel('K-band magnitude K')
188 frame1.legend(loc=0)
189
190 ###Plot no.2: K-band magnitude versus z via idealized K,z-relation and simulated data
191 around it.
192 fig2 = figure(figsize=(40/2.54, 40/2.54))
193 frame2 = fig2.add_subplot(1,1,1)
194 frame2.set_title("K-band magnitude vs z", fontsize=15)
195 frame2.plot(10*logzran ,Kran, 'g.', alpha=0.1, label="Simulated random distribution")
196 frame2.plot(10*logzarr ,Karr, linewidth=5, alpha=0.5, label="Idealized K,z-relation")
197 assymerr3 = [fdata5[5], fdata5[6]]
198 frame2.errorbar(fdata5[4], fdata5[0], yerr=fdata5[1], xerr=assymerr3, ecolor="k", elinewidth
199 =3, marker="*", mfc="y", ms=10, ls="") , label="Sample K, z")
200 frame2.set_xlabel('Redshift z')
201 frame2.set_ylabel('K-band magnitude K')
202 frame2.legend(loc=0)
203
204 ###Plot no.7: (data, 5 arcsec) log10(z) samples at some value of K and log10(z), showing
205 normal distributions.
206 fig7 = figure(figsize=(40/2.54, 40/2.54))
207 for i in range(len(fdata5[2])):
208     frame7 = fig7.add_subplot(4,4,i+1)
209     data = fdata5[7]["logzdist_{}".format(i)]
210     x = np.linspace(data.min(), data.max(), 100)
211     param = sps.norm.fit(data, floc=fdata5[2][i])
212     pdf_fitted = sps.norm.pdf(x, *param)
213     frame7.set_title("{}".format(newdatanames[i]), fontsize=22)
214     frame7.plot(x, pdf_fitted, color='m')
215     frame7.axvline(x=fdata5[2][i], linewidth=3, color='r') #####Mean
216     frame7.axvline(x=fdata5[2][i]-fdata5[3][i], linewidth=2, color='k') #####Std_logz
217     lower
218     frame7.axvline(x=fdata5[2][i]+fdata5[3][i], linewidth=2, color='k') #####Std_logz
219     upper
220     frame7.hist(data, normed=True, bins=50, alpha=0.3)
221     frame7.tick_params(labelsize=18)
222 fig7.tight_layout()
223
224 ###Plot no.8: (data, 5 arcsec) z samples at some value of K and z, showing lognormal
225 distributions.
226 fig8 = figure(figsize=(40/2.54, 40/2.54))
227 for i in range(len(fdata5[4])):
228     frame8 = fig8.add_subplot(4,4,i+1)
229     data = fdata5[8]["zdist_{}".format(i)]
230     x = np.linspace(data.min(), data.max(), 100)
231     param = sps.lognorm.fit(data, floc=0)
232     pdf_fitted = sps.lognorm.pdf(x, *param)
233     frame8.set_title("{}".format(newdatanames[i]), fontsize=22)
234     frame8.plot(x, pdf_fitted, color='m')
235     frame8.axvline(x=fdata5[4][i], linewidth=3, color='r') #####Median
236     frame8.axvline(x=fdata5[4][i]-fdata5[5][i], linewidth=2, color='k') #####Std_z lower
237     frame8.axvline(x=fdata5[4][i]+fdata5[6][i], linewidth=2, color='k') #####Std_z upper
238     frame8.hist(data, normed=True, bins=50, alpha=0.3)
239     frame8.tick_params(labelsize=18)
240 fig8.tight_layout()
241
242 show()

```

## B.4 Skymap creator

The following script is used to obtain skymaps of the source positions and predicted densities of GPS.

Scripts/Thesis\_synthesis3\_v2-4.py

```
1 #!/usr/bin/env python
2 from __future__ import division
3 import numpy as np
4 import numpy.random as npr
5 import scipy.optimize as spo
6 import fnmatch as fm
7 from matplotlib.pyplot import figure, show
8 from mpl_toolkits.mplot3d import Axes3D
9 #####THIS SCRIPT COMPUTES fun.
10
11
12 #####Synthesis I: Reading Data.
13 #####
14 print(">>> Synthesis I")
15
16 #####Reading Tjoa&McKean's name and data (because it's largely monocolumnar, this takes
17 some time and messy coding).
18 dat0TJOA = np.genfromtxt("poslist-McKean.dat", dtype=None, usecols=[0], comments="#")
19 dat1TJOA = np.genfromtxt("poslist-McKean.dat", dtype=None, usecols=[1], comments="#")
20 dat2TJOA = np.genfromtxt("poslist-McKean.dat", dtype=None, usecols=[2], comments="#")
21
22 namesTJOA, RAhrsTJOA, decdegTJOA = np.split(dat0TJOA, 3)
23 namesTJOA = namesTJOA.astype(str).flatten()
24 RAhrsTJOA = np.asarray(RAhrsTJOA.astype(float).flatten())
25 decdegTJOA = np.asarray(decdegTJOA.astype(float).flatten())
26
27 poep1TJOA, RAminTJOA, decminTJOA = np.split(dat1TJOA, 3)
28 RAminTJOA = np.asarray(RAminTJOA.astype(float).flatten())
29 decminTJOA = np.asarray(decminTJOA.astype(float).flatten())
30
31 poep2TJOA, RAsecTJOA, decsecTJOA = np.split(dat2TJOA, 3)
32 RAsecTJOA = np.asarray(RAsecTJOA.astype(float).flatten())
33 decsecTJOA = np.asarray(decsecTJOA.astype(float).flatten())
34
35 #####Reading Tjoa&McKean's synch and synch+powlaw data for spectral peaks.
36 purenamesTJOA = np.genfromtxt("ThesisTable_synchrotronNEW.dat", dtype=str, usecols=[1],
37 comments="#")
38 combinamesTJOA = np.genfromtxt("ThesisTable_synch-powlawNEW.dat", dtype=str, usecols=[1],
39 comments="#")
40 purevmaxTJOA = np.genfromtxt("ThesisTable_synchrotronNEW.dat", dtype=float, usecols=[10],
41 comments="#")
42 combivmaxTJOA = np.genfromtxt("ThesisTable_synch-powlawNEW.dat", dtype=float, usecols
43 =[14], comments="#")
44 vmxTJOA = np.concatenate((purevmaxTJOA, combivmaxTJOA))
45 pureSmaxTJOA = np.genfromtxt("ThesisTable_synchrotronNEW.dat", dtype=float, usecols=[12],
46 comments="#")
47 combiSmaxTJOA = np.genfromtxt("ThesisTable_synch-powlawNEW.dat", dtype=float, usecols
48 =[16], comments="#")
49 SmaxTJOA = np.concatenate((pureSmaxTJOA, combiSmaxTJOA))
50
51
52 #####Synthesis II: Reloaded. Computation of results.
53 #####
54 print(">>> Synthesis II: Reloaded")
55
56 def posmaker(RAhrs, RAmin, RAsec, decdeg, decmin, decsec): #####Function that computes the RA
57 and dec positions in fractional hours and degrees respectively for plotting purposes
58 RApos = RAhrs + RAmin/60 + (RAsec/60)/60
59 decpos = decdeg + decmin/60 + (decsec/60)/60
60 return RApos, decpos
61
62 def hidimpos(RA, dec): #####Function that projects RA and dec
63 data onto a sphere
64 x = np.cos(dec*np.pi/180)*np.sin(15*RA*np.pi/180)
65 y = np.cos(dec*np.pi/180)*np.cos(15*RA*np.pi/180)
```

```

55     z = np.sin(dec*np.pi/180)
56     return x,y,z
57
58 def areacalc(RAlow,RAupp,declow,decupp):
59     declow = declow*np.pi/180 ; decupp = decupp*np.pi/180
60     RAlow = 15*RAlow*np.pi/180 ; RAupp = 15*RAupp*np.pi/180
61     return (RAupp - RAlow)*(np.sin(decupp) - np.sin(declow))
62
63 ###Computing the positions of Tjoa&McKean's, Marlow's and Willott's sources.
64 posTJOA = posmaker(RAhrsTJOA,RAminTJOA,RAsecTJOA,decdegTJOA,decminTJOA,decsecTJOA)
65
66 ###Computing positions of pure and combined samples
67 pureRATJOA = []
68 puredecTJOA = []
69 for i in range(len(purenamesTJOA)):
70     for j in range(len(namesTJOA)):
71         if fm.fnmatch(purenamesTJOA[i],"{}".format(namesTJOA[j])) is True:
72             pureRATJOA.append(posTJOA[0][j])
73             puredecTJOA.append(posTJOA[1][j])
74 pureRATJOA = np.asarray(pureRATJOA)
75 puredecTJOA = np.asarray(puredecTJOA)
76
77 combiRATJOA = []
78 combidecTJOA = []
79 for i in range(len(combinamesTJOA)):
80     for j in range(len(namesTJOA)):
81         if fm.fnmatch(combinamesTJOA[i],"{}".format(namesTJOA[j])) is True:
82             combiRATJOA.append(posTJOA[0][j])
83             combidecTJOA.append(posTJOA[1][j])
84 combiRATJOA = np.asarray(combiRATJOA)
85 combidecTJOA = np.asarray(combidecTJOA)
86
87 peaknamesTJOA = list(purenamesTJOA)
88 peaknamesTJOA.extend(combinamesTJOA)
89 peakSmaxTJOA = np.concatenate((pureSmaxTJOA,combiSmaxTJOA))
90 peakRATJOA = np.concatenate((pureRATJOA,combiRATJOA))
91 peakdecTJOA = np.concatenate((puredecTJOA,combidecTJOA))
92
93 ###Computing Mirach and Grumium's positions as reference material for skymaps.
94 Mirachpos = posmaker(1.0,9.0,43.92388,35.0,37.0,14.0075)
95 Grumiumpos = posmaker(17.0,53.0,31.72962,56.0,52.0,21.5143)
96
97 ###Loop that splits Tjoa&McKean's data (all of it) into its two component fields of view
98     (1<RA<2, 35<dec<45 low and 17<RA<18, 55<dec<60 high).
99 RAhigh = [] ; dechigh = []
100 RAlow = [] ; declow = []
101 for i in range(len(posTJOA[0])):
102     if posTJOA[0][i] > 2.0:
103         RAhigh.append(posTJOA[0][i])
104         dechigh.append(posTJOA[1][i])
105     elif posTJOA[0][i] < 2.0:
106         RAlow.append(posTJOA[0][i])
107         declow.append(posTJOA[1][i])
108
109 ###Loop that splits Tjoa&McKean's data (only peaked) into its two component fields of
110     view (1<RA<2, 35<dec<45 low and 17<RA<18, 55<dec<60 high).
111 RAhighp = [] ; dechighp = [] ; Smaxhigh = []
112 RAlowp = [] ; declowp = [] ; Smaxlow = []
113 for i in range(len(peakRATJOA)):
114     if peakRATJOA[i] > 2.0:
115         RAhighp.append(peakRATJOA[i])
116         dechighp.append(peakdecTJOA[i])
117         Smaxhigh.append(peakSmaxTJOA[i])
118     elif peakRATJOA[i] < 2.0:
119         RAlowp.append(peakRATJOA[i])
120         declowp.append(peakdecTJOA[i])
121         Smaxlow.append(peakSmaxTJOA[i])
122
123 def intnumdensflux(Slow,Sp):
124     k = 0

```

```

123     for i in range(len(Sp)):
124         if Sp[i] >= Slow:
125             k = k + 1
126         else:
127             print("denied")
128     return k
129
130 def difnumdensflux(Slow, Supp, Sp):
131     k = 0
132     for i in range(len(Sp)):
133         if Sp[i] >= Slow and Sp[i] <= Supp:
134             k = k + 1
135         else:
136             print("denied")
137     return k
138
139 ###Computing densities for both fields and subsequently computing the total estimated
140     amount of sources of this calibre.
141 RAlowhigh = 17 ; RAupphigh = 18 ; declowhigh = 55 ; decupphigh = 60
142 RAlowlow = 1 ; RAupplow = 2 ; declowlow = 35 ; decupplow = 45
143 areahigh = areacalc(RAlowhigh, RAupphigh, declowhigh, decupphigh)
144 arealow = areacalc(RAlowlow, RAupplow, declowlow, decupplow)
145 denshigh = len(RAhigh)/areahigh
146 denslow = len(RAlow)/arealow
147 densav = (denshigh+denslow)/2
148 totalsources = 4*np.pi*densav
149
150 print("Area of high RA field is:", areahigh, "steradians")
151 print("Area of low RA field is:", arealow, "steradians")
152 print("Source density in high RA field is:", denshigh, "per steradian")
153 print("Source density in low RA field is:", denslow, "per steradian")
154 print("Source density average is:", densav, "per steradian")
155 print("So total amount of observable sources in the sky with flux at 5 GHz of ~50 mJy
156     should be:", totalsources)
157
158 ###Computing number density per steradian as a function of peak flux (integrated method)
159 Slow1 = np.asarray([20.0, 40.0, 60.0, 80.0])
160 nsrinthigh = []
161 nsrintlow = []
162 for i in range(len(Slow1)):
163     x1 = intnumdensflux(Slow1[i], Smaxhigh)
164     x2 = intnumdensflux(Slow1[i], Smaxlow)
165     nsrinthigh.append(x1)
166     nsrintlow.append(x2)
167 nsrinthigh = np.asarray(nsrinhigh)
168 nsrintlow = np.asarray(nsrintlow)
169 densSinhigh = nsrinhigh/areahigh
170 densSintlow = nsrintlow/arealow
171 densSintav = (densSinhigh+densSintlow)/2
172 totalSintsources = 4*np.pi*densSintav
173 projSintsources = [totalSintsources[0]-totalSintsources[1], totalSintsources[1]-
174     totalSintsources[2], totalSintsources[2]-totalSintsources[3], totalSintsources[3]]
175
176 ###Computing number density-peak flux convolution per steradian as a function of peak
177     flux (differential method).
178 Smed = np.asarray([30.0, 60.0])
179 Sbinsize = np.asarray([10.0, 20.0])
180 Slow = Smed-Sbinsize
181 Supp = Smed+Sbinsize
182 nsrdifhigh = []
183 nsrdiflow = []
184 for i in range(len(Smed)):
185     x1 = difnumdensflux(Slow[i], Supp[i], Smaxhigh)
186     x2 = difnumdensflux(Slow[i], Supp[i], Smaxlow)
187     nsrdifhigh.append(x1)
188     nsrdiflow.append(x2)
189 nsrdifhigh = np.asarray(nsrdifhigh)
190 nsrdiflow = np.asarray(nsrdiflow)

```

```

188 nsrdifav = nsrdifhigh+nsrdiflow
189 densSdifhigh = nsrdifhigh/areahigh
190 densSdiflow = nsrdiflow/arealow
191 densSdifav = nsrdifav/(areahigh+arealow)
192 projSdifsources = 4*np.pi*densSdifav
193
194 print (nsrdiflow)
195 print (nsrdifhigh)
196
197 ###Snellen data.
198 SmedSNEL = np.asarray([75,150,300])
199 plotSpeakSNEL = np.asarray([1.05089459,2.37790302,5.88500972])
200 SbinsizeSNEL = np.asarray([25.0,50.0,100.0])
201 plotSerrSNEL = np.asarray([0.33232205,0.84071567,2.2243246])
202
203 ###AVENSS data.
204 SmedVRIES = np.asarray([4000])
205 plotSpeakVRIES = np.asarray([22.0])
206 SbinsizeVRIES = np.asarray([2000.0])
207 plotSerrVRIES = np.asarray([10.0])
208
209 ###Computing 3d positions on sky for star orb.
210 p3dTJOA = hidimpos(-posTJOA[0],posTJOA[1])
211
212 ###Computing 3d positions of predicted sources (locations are not actually physical but
the density is).
213 Phi = 2*np.pi*npr.random(size=totalsources)/15*180/np.pi
214 Theta = np.arccos(2*npr.random(size=totalsources)-1)*180/np.pi-90
215 p3dPRED = hidimpos(-Phi,Theta)
216
217 Phi0 = 2*np.pi*npr.random(size=totalsources-totalSintsources[0])/15*180/np.pi
218 Theta0 = np.arccos(2*npr.random(size=totalsources-totalSintsources[0])-1)*180/np.pi-90
219 p3dPRED0 = hidimpos(-Phi0,Theta0)
220
221 Phi2 = 2*np.pi*npr.random(size=projSintsources[0])/15*180/np.pi
222 Theta2 = np.arccos(2*npr.random(size=projSintsources[0])-1)*180/np.pi-90
223 p3dPRED2 = hidimpos(-Phi2,Theta2)
224
225 Phi4 = 2*np.pi*npr.random(size=projSintsources[1])/15*180/np.pi
226 Theta4 = np.arccos(2*npr.random(size=projSintsources[1])-1)*180/np.pi-90
227 p3dPRED4 = hidimpos(-Phi4,Theta4)
228
229 Phi6 = 2*np.pi*npr.random(size=projSintsources[2])/15*180/np.pi
230 Theta6 = np.arccos(2*npr.random(size=projSintsources[2])-1)*180/np.pi-90
231 p3dPRED6 = hidimpos(-Phi6,Theta6)
232
233 Phi8 = 2*np.pi*npr.random(size=projSintsources[3])/15*180/np.pi
234 Theta8 = np.arccos(2*npr.random(size=projSintsources[3])-1)*180/np.pi-90
235 p3dPRED8 = hidimpos(-Phi8,Theta8)
236
237 dndS = densSdifav/(Supp/1000-Slow/1000)
238 plotSpeak = dndS*(Smed/1000)**(5/2)
239 plotSerr = (np.sqrt(nsrdifav)/nsrdifav)*plotSpeak
240 print (plotSpeak)
241 print (plotSerr)
242
243 def fitfun(S,a,b,c):
244     x = np.log10(S)
245     return 10**(a + b*x + c*x**2)
246
247 Sarr = np.arange(10,10000,0.1)
248 Slst = np.concatenate((Smed,SmedSNEL,SmedVRIES))
249 peaklst = np.concatenate((plotSpeak,plotSpeakSNEL,plotSpeakVRIES))
250 peakerr = np.concatenate((plotSerr,plotSerrSNEL,plotSerrVRIES))
251 p,pcov = spo.curve_fit(fitfun,Slst,peaklst,sigma=peakerr,**{'maxfev':1000000})
252 print (p)
253 print (pcov)
254 function = fitfun(Sarr,*p)
255
256 #####Synthesis III: Revolutions. Plotting the results.

```

```

#####
257 print(">>> Synthesis III: Revolutions")
258
259 ###Plot no.1: number of sources per area.
260 fig1 = figure(figsize=(40/2.54, 40/2.54))
261 frame1 = fig1.add_subplot(1,1,1)
262 frame1.set_title("Source position in sky",fontsize=15)
263 frame1.set_xlim([RALowlow,RAUpplhigh])
264 frame1.set_ylim([declowlow,decupplhigh])
265 frame1.axvline(x=RALowlow,linewidth=1,color='b') ; frame1.axvline(x=RAUpplow,linewidth
    =1,color='b')
266 frame1.axhline(y=declowlow,linewidth=1,color='b') ; frame1.axhline(y=decupplow,linewidth
    =1,color='b')
267 frame1.axvline(x=RAlowhigh,linewidth=1,color='r') ; frame1.axvline(x=RAUpplhigh,linewidth
    =1,color='r')
268 frame1.axhline(y=declowhigh,linewidth=1,color='r') ; frame1.axhline(y=decupplhigh,
    linewidth=1,color='r')
269 frame1.plot(posTJOA[0],posTJOA[1], 'y*',ms=11)
270 frame1.set_xlabel('Right Ascension (hours)')
271 frame1.set_ylabel('Declination (deg)')
272 frame1.invert_xaxis()
273
274 ###Plot no.2: number of sources per area, closeup low RA.
275 fig2 = figure(figsize=(40/2.54, 40/2.54))
276 frame2 = fig2.add_subplot(1,1,1)
277 frame2.set_title("Source position in sky",fontsize=15)
278 frame2.set_xlim([RALowlow,RAUpplow])
279 frame2.set_ylim([declowlow,decupplow])
280 frame2.plot(posTJOA[0],posTJOA[1], 'y*',ms=15)
281 frame2.plot(peakRATJOA,peakdecTJOA, 'c*',ms=15)
282 frame2.plot(Mirachpos[0],Mirachpos[1], 'r*',ms=20)
283 frame2.annotate("Mirach (Beta Andromedae)",xy=(Mirachpos[0]+0.05,Mirachpos[1]+0.1))
284 frame2.set_xlabel('Right Ascension (hours)')
285 frame2.set_ylabel('Declination (deg)')
286 frame2.invert_xaxis()
287
288 ###Plot no.3: number of sources per area, closeup high RA.
289 fig3 = figure(figsize=(40/2.54, 40/2.54))
290 frame3 = fig3.add_subplot(1,1,1)
291 frame3.set_title("Source position in sky",fontsize=15)
292 frame3.set_xlim([RAlowhigh,RAUpplhigh])
293 frame3.set_ylim([declowhigh,decupplhigh])
294 frame3.plot(posTJOA[0],posTJOA[1], 'y*',ms=15)
295 frame3.plot(peakRATJOA,peakdecTJOA, 'c*',ms=15)
296 frame3.plot(Grumiumpos[0],Grumiumpos[1], 'b*',ms=20)
297 frame3.annotate("Grumium (Xi Draconis)",xy=(Grumiumpos[0]+0.05,Grumiumpos[1]-0.2))
298 frame3.set_xlabel('Right Ascension (hours)')
299 frame3.set_ylabel('Declination (deg)')
300 frame3.invert_xaxis()
301
302 ###Plot no.4: sources both real and predicted projected on sphere, assuming density
    found from McKean data.
303 fig4 = figure(figsize=(40/2.54, 40/2.54))
304 frame4 = fig4.add_subplot(111, projection='3d')
305 frame4.plot(p3dPRED[0],p3dPRED[1],p3dPRED[2], 'y*',markersize=8)
306 frame4.plot(p3dTJOA[0],p3dTJOA[1],p3dTJOA[2], 'r*',markersize=12)
307 frame4.set_title('Isotropic predicted spherical distribution of GPS sources')
308 frame4.set_xlabel("x")
309 frame4.set_ylabel("y")
310 frame4.set_zlabel("z")
311
312 ###Plot no.5: sources both real and predicted projected on plane.
313 fig5 = figure(figsize=(40/2.54, 40/2.54))
314 frame5 = fig5.add_subplot(1, 1, 1)
315 frame5.plot(Phi,Theta, 'y*',markersize=8,label="Predicted source density")
316 frame5.plot(posTJOA[0],posTJOA[1], 'r*',markersize=12,label="Our data")
317 frame5.set_title('Isotropic predicted distribution of GPS sources over flattened sky')
318 frame5.set_xlabel("Right Ascension (hours)")
319 frame5.set_ylabel("Declination (degrees)")
320 frame5.invert_xaxis()

```

```

321
322 ###Plot no.6: number density as a function of peak flux.
323 fig6 = figure(figsize=(40/2.54, 40/2.54))
324 frame6 = fig6.add_subplot(1, 1, 1)
325 frame6.set_xlim([0,100])
326 frame6.plot(Slow1, densSintav, 'r*', markersize=15, label="averaged sample")
327 frame6.set_title('cumulative number density as a function of peak flux')
328 frame6.set_xlabel("Smax (mJy)")
329 frame6.set_ylabel("number density with Smax in excess of S-value")
330 frame6.legend()
331
332 ###Plot no.7: number as a function of peak flux.
333 fig7 = figure(figsize=(40/2.54, 40/2.54))
334 frame7 = fig7.add_subplot(1, 1, 1)
335 frame7.set_xlim([0,100])
336 frame7.plot(Slow1, totalSintsources, 'r*', markersize=15, label="predicted sources")
337 frame7.set_title('predicted cumulative number as a function of peak flux')
338 frame7.set_xlabel("Smax (mJy)")
339 frame7.set_ylabel("number with Smax in excess of S-value")
340 frame7.legend()
341
342 ###Plot no.8: predicted sources separated by peak flux population projected on sphere,
      assuming density found from McKean data.
343 fig8 = figure(figsize=(40/2.54, 40/2.54))
344 frame8 = fig8.add_subplot(111, projection='3d')
345 #frame8.plot(p3dPRED0[0], p3dPRED0[1], p3dPRED0[2], 'm*', markeredgewidth=0, markersize=5,
      label='No peak', alpha=0.5)
346 frame8.plot(p3dPRED2[0], p3dPRED2[1], p3dPRED2[2], 'r*', markeredgewidth=0, markersize=7,
      label='20-40 mJy')
347 frame8.plot(p3dPRED4[0], p3dPRED4[1], p3dPRED4[2], 'y*', markeredgewidth=0, markersize=7,
      label='40-60 mJy')
348 frame8.plot(p3dPRED6[0], p3dPRED6[1], p3dPRED6[2], 'c*', markeredgewidth=0, markersize=7,
      label='60-80 mJy')
349 frame8.plot(p3dPRED8[0], p3dPRED8[1], p3dPRED8[2], 'b*', markeredgewidth=0, markersize=7,
      label='>80 mJy')
350 frame8.set_title('Isotropic predicted spherical distribution of GPS sources')
351 frame8.set_xlabel("x")
352 frame8.set_ylabel("y")
353 frame8.set_zlabel("z")
354 frame8.legend()
355
356 ###Plot no.9: predicted sources projected on plane.
357 fig9 = figure(figsize=(40/2.54, 40/2.54))
358 frame9 = fig9.add_subplot(1, 1, 1)
359 #frame9.plot(Phi0, Theta0, 'm*', markeredgewidth=0, markersize=5, label='No peak', alpha=0.5)
360 frame9.plot(Phi2, Theta2, 'r*', markeredgewidth=0, markersize=7, label='20-40 mJy')
361 frame9.plot(Phi4, Theta4, 'y*', markeredgewidth=0, markersize=7, label='40-60 mJy')
362 frame9.plot(Phi6, Theta6, 'c*', markeredgewidth=0, markersize=7, label='60-80 mJy')
363 frame9.plot(Phi8, Theta8, 'b*', markeredgewidth=0, markersize=7, label='>80 mJy')
364 frame9.set_title('Isotropic predicted distribution of GPS sources over flattened sky')
365 frame9.set_xlabel("Right Ascension (hours)")
366 frame9.set_ylabel("Declination (degrees)")
367 frame9.legend()
368 frame9.invert_xaxis()
369
370
371 show()

```

## B.5 Number count calculator

The following script is used to obtain number counts.

Scripts/Thesis\_synthesis4\_v1-2.py

```

1 #!/usr/bin/env python
2 from __future__ import division
3 import numpy as np
4 import scipy as sp
5 import scipy.stats as sps

```

```

6 import fmatch as fm
7 import numpy.random as npr
8 from matplotlib.pyplot import figure, show
9 ###THIS SCRIPT COMPUTES STD IN LOG10(z) AND z.
10
11
12 #####Synthesis I: Reading Data.
13 #####
14 print(">>> Synthesis I")
15
16 ###McKean's & Tjoa's data (K,z)
17 znamesTJOA = np.genfromtxt("K,z-data-fulltable.dat", dtype=str, usecols=[0], comments="#")
18 zdatTJOA = np.genfromtxt("K,z-data-fulltable.dat", usecols=[1], comments="#")
19 zstdlowTJOA = np.genfromtxt("K,z-data-fulltable.dat", usecols=[2], comments="#")
20 zstduppTJOA = np.genfromtxt("K,z-data-fulltable.dat", usecols=[3], comments="#")
21 KdatTJOA = np.genfromtxt("K,z-data-fulltable.dat", usecols=[4], comments="#")
22 KstdTJOA = np.genfromtxt("K,z-data-fulltable.dat", usecols=[5], comments="#")
23
24 purenamesTJOA = np.genfromtxt("ThesisTable_synchrotronNEW.dat", dtype=str, usecols=[1],
25     comments="#")
26 combinamesTJOA = np.genfromtxt("ThesisTable_synch-powlawNEW.dat", dtype=str, usecols=[1],
27     comments="#")
28 purevmaxTJOA = np.genfromtxt("ThesisTable_synchrotronNEW.dat", dtype=float, usecols=[10],
29     comments="#")
30 combivmaxTJOA = np.genfromtxt("ThesisTable_synch-powlawNEW.dat", dtype=float, usecols
31     =[14], comments="#")
32 vmaxTJOA = np.concatenate((purevmaxTJOA, combivmaxTJOA))
33 pureSmaxTJOA = np.genfromtxt("ThesisTable_synchrotronNEW.dat", dtype=float, usecols=[12],
34     comments="#")
35 combiSmaxTJOA = np.genfromtxt("ThesisTable_synch-powlawNEW.dat", dtype=float, usecols
36     =[16], comments="#")
37 SmaxTJOA = np.concatenate((pureSmaxTJOA, combiSmaxTJOA))
38
39 #####Synthesis II: Reloaded. Computation of results.
40 #####
41 print(">>> Synthesis II: Reloaded")
42
43 def Kzrel(z):
44     return 17.37 + 4.53*np.log10(z) + 0.31*(np.log10(z))**2
45
46 ###Generating plain, idealized K-log10(z) relation.
47 zarr = np.linspace(0.05,5,10000)
48 Karr = Kzrel(zarr)
49
50 znamesTJOA2 = []
51 zdatTJOA2 = []
52 zstdlowTJOA2 = []
53 zstduppTJOA2 = []
54 KdatTJOA2 = []
55 KstdTJOA2 = []
56 for i in range(len(zdatTJOA)):
57     if zdatTJOA[i] == 0.0:
58         print("This point has a nonexistent redshift.")
59     else:
60         znamesTJOA2.append(znamesTJOA[i])
61         zdatTJOA2.append(zdatTJOA[i])
62         zstdlowTJOA2.append(zstdlowTJOA[i])
63         zstduppTJOA2.append(zstduppTJOA[i])
64         KdatTJOA2.append(KdatTJOA[i])
65         KstdTJOA2.append(KstdTJOA[i])
66 znamesTJOA2 = np.asarray(znamesTJOA2)
67 zdatTJOA2 = np.asarray(zdatTJOA2)
68 zstdlowTJOA2 = np.asarray(zstdlowTJOA2)
69 zstduppTJOA2 = np.asarray(zstduppTJOA2)
70 KdatTJOA2 = np.asarray(KdatTJOA2)
71 KstdTJOA2 = np.asarray(KstdTJOA2)
72
73 znamesTJOApure = []
74 zdatTJOApure = []

```

```

68 zstdlowTJOApure = []
69 zstduppTJOApure = []
70 KdatTJOApure = []
71 KstdTJOApure = []
72 for i in range(len(purenamesTJOA)):
73     for j in range(len(znamesTJOA2)):
74         if fm.fnmatch(purenamesTJOA[i], "{}".format(znamesTJOA2[j])) is True: ###Check
75             whether the chosen name in z-data matches any known pure spectral peak name
76             znamesTJOApure.append(znamesTJOA2[j]) ###If so, append
77             indices to specnames- and Knames lists.
78             zdatTJOApure.append(zdatTJOA2[j])
79             zstdlowTJOApure.append(zstdlowTJOA2[j])
80             zstduppTJOApure.append(zstduppTJOA2[j])
81             KdatTJOApure.append(KdatTJOA2[j])
82             KstdTJOApure.append(KstdTJOA2[j])
83 znamesTJOApure = np.asarray(znamesTJOApure)
84 zdatTJOApure = np.asarray(zdatTJOApure)
85 zstdlowTJOApure = np.asarray(zstdlowTJOApure)
86 zstduppTJOApure = np.asarray(zstduppTJOApure)
87 KdatTJOApure = np.asarray(KdatTJOApure)
88 KstdTJOApure = np.asarray(KstdTJOApure)
89
90 znamesTJOAcombi = []
91 zdatTJOAcombi = []
92 zstdlowTJOAcombi = []
93 zstduppTJOAcombi = []
94 KdatTJOAcombi = []
95 KstdTJOAcombi = []
96 for i in range(len(combinamesTJOA)):
97     for j in range(len(znamesTJOA2)):
98         if fm.fnmatch(combinamesTJOA[i], "{}".format(znamesTJOA2[j])) is True: ###Check
99             whether the chosen name in z-data matches any known pure spectral peak name
100             znamesTJOAcombi.append(znamesTJOA2[j]) ###If so, append
101             indices to specnames- and Knames lists.
102             zdatTJOAcombi.append(zdatTJOA2[j])
103             zstdlowTJOAcombi.append(zstdlowTJOA2[j])
104             zstduppTJOAcombi.append(zstduppTJOA2[j])
105             KdatTJOAcombi.append(KdatTJOA2[j])
106             KstdTJOAcombi.append(KstdTJOA2[j])
107 znamesTJOAcombi = np.asarray(znamesTJOAcombi)
108 zdatTJOAcombi = np.asarray(zdatTJOAcombi)
109 zstdlowTJOAcombi = np.asarray(zstdlowTJOAcombi)
110 zstduppTJOAcombi = np.asarray(zstduppTJOAcombi)
111 KdatTJOAcombi = np.asarray(KdatTJOAcombi)
112 KstdTJOAcombi = np.asarray(KstdTJOAcombi)
113
114 #####Synthesis III: Revolutions. Plotting the results.
115 #####
116 print(">>> Synthesis III: Revolutions")
117
118 ###Plot no.1: K-band magnitude versus z.
119 fig1 = figure(figsize=(40/2.54, 40/2.54))
120 frame1 = fig1.add_subplot(1,1,1)
121 frame1.set_title("K-band magnitude vs z", fontsize=20)
122 frame1.plot(zarr, Karr, "k", label="Idealized K,z-relation")
123 assymerrzTJOA2 = [zstdlowTJOA2, zstduppTJOA2]
124 frame1.errorbar(zdatTJOA2, KdatTJOA2, xerr=assymerrzTJOA2, yerr=KstdTJOA2, ecolor="k",
125               elinewidth=1, marker="*", mfc="y", ms=12, ls="", label="Our data points (K,z)")
126 frame1.set_xlabel('Redshift z', fontsize=20)
127 frame1.set_ylabel('K-band magnitude', fontsize=20)
128 frame1.tick_params(labelsize=20)
129
130 ###Plot no.2: number of sources per redshift bin (pure and combined separately).
131 fig2 = figure(figsize=(40/2.54, 40/2.54))
132 frame2 = fig2.add_subplot(1,1,1)
133 frame2.set_title("number of sources per $z$", fontsize=20)
134 frame2.hist(zdatTJOApure, bins=[0.0,0.5,1.0,1.5,2.0,2.5,3.0], label='GPS', alpha=0.5)
135 frame2.hist(zdatTJOAcombi, bins=[0.0,0.5,1.0,1.5,2.0,2.5,3.0], label='GPS + power law',
136           alpha=0.5)

```

```

131 #for i in zdatTJOApure:
132 #     frame2.axvline(x=i,linewidth=2,color='y')
133 #for i in zdatTJOAcombi:
134 #     frame2.axvline(x=i,linewidth=2,color='g')
135 frame2.set_xlabel('Redshift $z$', fontsize=20)
136 frame2.set_ylabel('Number of sources', fontsize=20)
137 frame2.tick_params(labelsize=20)
138
139 #####Plot no.3: number of sources per redshift bin (all of them).
140 fig3 = figure(figsize=(40/2.54, 40/2.54))
141 frame3 = fig3.add_subplot(1,1,1)
142 frame3.set_title("number of sources per $z$", fontsize=20)
143 frame3.hist(zdatTJOA2, bins=[0.0,0.5,1.0,1.5,2.0,2.5,3.0])
144 #for i in zdatTJOA2:
145 #     frame3.axvline(x=i,linewidth=1,color='r')
146 frame3.set_xlabel('Redshift $z$', fontsize=20)
147 frame3.set_ylabel('Number of sources', fontsize=20)
148 frame3.tick_params(labelsize=20)
149
150 #####Plot no.4: number of sources per redshift bin (pure and combined separately).
151 fig4 = figure(figsize=(40/2.54, 40/2.54))
152 frame4 = fig4.add_subplot(1,1,1)
153 frame4.set_title("number of sources per $z$", fontsize=20)
154 frame4.hist(np.log10(zdatTJOApure), bins=[-2.0,-1.5,-1.0,-0.5,0.0,0.5,1.0,1.5,2.0], label=
155             'GPS', alpha=0.5)
156 frame4.hist(np.log10(zdatTJOAcombi), bins=[-2.0,-1.5,-1.0,-0.5,0.0,0.5,1.0,1.5,2.0], label=
157             'GPS + power law', alpha=0.5)
158 #for i in zdatTJOApure:
159 #     frame4.axvline(x=np.log10(i),linewidth=1,color='y')
160 #for i in zdatTJOAcombi:
161 #     frame4.axvline(x=np.log10(i),linewidth=1,color='g')
162 frame4.set_xlabel('Redshift log10z', fontsize=20)
163 frame4.set_ylabel('Number of sources', fontsize=20)
164 frame4.tick_params(labelsize=20)
165
166 #####Plot no.5: number of sources per redshift bin (all of them).
167 fig5 = figure(figsize=(40/2.54, 40/2.54))
168 frame5 = fig5.add_subplot(1,1,1)
169 frame5.set_title("number of sources per $z$", fontsize=20)
170 frame5.hist(np.log10(zdatTJOA2), bins=[-2.0,-1.5,-1.0,-0.5,0.0,0.5,1.0,1.5,2.0])
171 #for i in zdatTJOA2:
172 #     frame5.axvline(x=np.log10(i),linewidth=1,color='r')
173 frame5.set_xlabel('Redshift log10z', fontsize=20)
174 frame5.set_ylabel('Number of sources', fontsize=20)
175 frame5.tick_params(labelsize=20)
176
177 #####Plot no.6: number of sources per vmax bin (pure and combined separately).
178 fig6 = figure(figsize=(40/2.54, 40/2.54))
179 frame6 = fig6.add_subplot(1,1,1)
180 frame6.set_title("number of sources per $v_{max}$", fontsize=20)
181 frame6.hist(purevmaxTJOA, bins=[0.0,2.0,4.0,6.0,8.0,10.0,12.0,14.0,16.0,18.0,20.0], label=
182             'GPS', alpha=0.5)
183 frame6.hist(combivmaxTJOA, bins=[0.0,2.0,4.0,6.0,8.0,10.0,12.0,14.0,16.0,18.0,20.0], label=
184             'GPS + power law', alpha=0.5)
185 #for i in purevmaxTJOA:
186 #     frame6.axvline(x=i,linewidth=2,color='y')
187 #for i in combivmaxTJOA:
188 #     frame6.axvline(x=i,linewidth=2,color='g')
189 frame6.set_xlabel('Peak frequency $v_{max}$ (GHz)', fontsize=20)
190 frame6.set_ylabel('Number of sources', fontsize=20)
191 frame6.tick_params(labelsize=20)
192
193 #####Plot no.7: number of sources per vmax bin (all of them).
194 fig7 = figure(figsize=(40/2.54, 40/2.54))
195 frame7 = fig7.add_subplot(1,1,1)
196 frame7.set_title("number of sources per $v_{max}$", fontsize=20)
197 frame7.hist(vmaxTJOA, bins=[0.0,2.0,4.0,6.0,8.0,10.0,12.0,14.0,16.0,18.0,20.0])
198 #for i in vmaxTJOA:
199 #     frame7.axvline(x=i,linewidth=2,color='r')
200 frame7.set_xlabel('Peak frequency $v_{max}$ (GHz)', fontsize=20)

```

```

197 frame7.set_ylabel('Number of sources', fontsize=20)
198 frame7.tick_params(labelsize=20)
199
200 ###Plot no.8: number of sources per Smax bin (pure and combined separately).
201 fig8 = figure(figsize=(40/2.54, 40/2.54))
202 frame8 = fig8.add_subplot(1,1,1)
203 frame8.set_title("number of sources per Smax", fontsize=20)
204 frame8.hist(pureSmaxTJOA, bins=[20.0,30.0,40.0,50.0,60.0,70.0,80.0,90.0], label='GPS',
205            alpha=0.5)
206 frame8.hist(combiSmaxTJOA, bins=[20.0,30.0,40.0,50.0,60.0,70.0,80.0,90.0], label='GPS +
207            power law', alpha=0.5)
208 #for i in pureSmaxTJOA:
209 #    frame8.axvline(x=i, linewidth=2, color='y')
210 #for i in combiSmaxTJOA:
211 #    frame8.axvline(x=i, linewidth=2, color='g')
212 frame8.set_xlabel('Peak flux Smax (mJy)', fontsize=20)
213 frame8.set_ylabel('Number of sources', fontsize=20)
214 frame8.tick_params(labelsize=20)
215
216 ###Plot no.9: number of sources per Smax bin (all of them).
217 fig9 = figure(figsize=(40/2.54, 40/2.54))
218 frame9 = fig9.add_subplot(1,1,1)
219 frame9.set_title("number of sources per Smax", fontsize=20)
220 frame9.hist(SmaxTJOA, bins=[20.0,30.0,40.0,50.0,60.0,70.0,80.0,90.0])
221 #for i in SmaxTJOA:
222 #    frame9.axvline(x=i, linewidth=2, color='r')
223 frame9.set_xlabel('Peak flux Smax (mJy)', fontsize=20)
224 frame9.set_ylabel('Number of sources', fontsize=20)
225 frame9.tick_params(labelsize=20)
226
227 show()

```

## B.6 Size & distance cosmological calculator

The following script is used to synthesize results to obtain luminosity distances and angular & linear sizes.

Scripts/Thesis\_synthesis6\_v2-3.py

```

1  #!/usr/bin/env python
2  from __future__ import division
3  import numpy as np
4  import scipy.optimize as spo
5  import fnmatch as fm
6  from cosmodistfun import cosmocalc
7  from matplotlib.pyplot import figure, show
8  ###THIS SCRIPT COMPUTES STD IN LOG10(z) AND z.
9
10
11 #####Synthesis I: Reading Data.
12 #####
13 print(">>> Synthesis I")
14
15 ###Reading relevant data from datafiles: McKean&Tjoa data.
16 znames = np.genfromtxt("K,z-data-ztable.dat", dtype=str, usecols=[0], comments="#")
17 zdat, zstdlow, zstdupp, Kdat, Kstd = np.genfromtxt("K,z-data-ztable.dat", usecols
18           =[1,2,3,4,5], unpack=True, comments="#")
19
20 spec1names = np.genfromtxt("ThesisTable_synchrotronNEW.dat", dtype=str, usecols=[1],
21           comments="#")
22 vmax1dat, vmax1std, Smax1dat, Smax1std = np.genfromtxt("ThesisTable_synchrotronNEW.dat",
23           dtype=float, usecols=[10,11,12,13], unpack=True, comments="#")
24 spec2names = np.genfromtxt("ThesisTable_synch-powlawNEW.dat", dtype=str, usecols=[1],
25           comments="#")
26 vmax2dat, vmax2std, Smax2dat, Smax2std = np.genfromtxt("ThesisTable_synch-powlawNEW.dat",
27           dtype=float, usecols=[14,15,16,17], unpack=True, comments="#")
28
29 ###Reading relevant data from datafiles: Snellen faint dataset.

```

```

24 specnamesSNELf = np.genfromtxt("Snellensample-faint.dat", dtype=str, usecols=[0], comments=
    "#")
25 zdatSNELf, vmaxdatSNELf, SmaxdatSNELf, angmarSNELf = np.genfromtxt("Snellensample-faint.dat",
    dtype=float, usecols=[1, 2, 3, 5], unpack=True, comments="#")
26
27 ###Reading relevant data from datafiles: Snellen bright dataset.
28 specnamesSNELb = np.genfromtxt("Snellensample-bright.dat", dtype=str, usecols=[0], comments
    ="#")
29 zdatSNELb, vmaxdatSNELb, SmaxdatSNELb, angmarSNELb = np.genfromtxt("Snellensample-bright.
    dat", dtype=float, usecols=[1, 2, 3, 5], unpack=True, comments="#")
30 SmaxdatSNELb = SmaxdatSNELb*1000
31
32 ###Reading relevant data from datafiles: Snellen CSS dataset.
33 specnamesSNELc = np.genfromtxt("Snellensample-CSS.dat", dtype=str, usecols=[0], comments="#
    ")
34 zdatSNELc, vmaxdatSNELc, SmaxdatSNELc, angmarSNELc = np.genfromtxt("Snellensample-CSS.dat",
    dtype=float, usecols=[1, 2, 3, 5], unpack=True, comments="#")
35 SmaxdatSNELc = SmaxdatSNELc*1000
36 angmarSNELc = angmarSNELc*1000
37
38 ###Concatenating arrays.
39 typelst = []
40 for i in range(len(vmax1dat)):
41     typelst.append("GPS")
42 for i in range(len(vmax2dat)):
43     typelst.append("GPPL")
44
45 vmaxdat = np.concatenate((vmax1dat, vmax2dat)) ; vmaxstd = np.concatenate((vmax1std,
    vmax2std))                                     ###Concatenating
    vmax data and std's
46 Smaxdat = np.concatenate((Smax1dat, Smax2dat)) ; Smaxstd = np.concatenate((Smax1std,
    Smax2std))                                     ###Concatenating
    Smax data and std's
47
48 specnames = []                                     ###Creating spectral names array
49 specnames.extend(spec1names)                       ###Appending pure synch names
50 specnames.extend(spec2names)                       ###Appending synch+powlaw names
51
52 ###Loop to create an array with the peak data for those sources whose redshift is known:
    McKean&Tjoa data.
53 names = []                                         ###Empty names list
54 Smaxlst = []                                       ###Empty S-peak list
55 vmaxlst = []                                       ###Empty v-peak list
56 Smaxstdlst = []                                    ###Empty S-peak std list
57 vmaxstdlst = []                                    ###Empty v-peak std list
58 zlst = []                                          ###Empty z list
59 zstdlowlst = [] ; zstdupplst = []                 ###Empty z lower- and upper boundaries lists
60 for i in range(len(specnames)):
61     k = 0
62     names.append(specnames[i])                     ###Appending name to name list
63     Smaxlst.append(Smaxdat[i])                     ###Appending relevant entries from
    Smax-data
64     vmaxlst.append(vmaxdat[i])                     ###Appending relevant entries from
    vmax-data
65     Smaxstdlst.append(Smaxstd[i])                  ###Appending relevant entries from Smax-
    std-data
66     vmaxstdlst.append(vmaxstd[i])                  ###Appending relevant entries from vmax-
    std-data
67     for j in range(len(znames)):
68         if fm.fnmatch(specnames[i], "{}".format(znames[j])) is True:             ###Check
    whether the chosen name in z-data matches any known spectral peak name
69             zlst.append(zdat[j])                 ###Appending relevant
    entries from z-data
70             zstdlowlst.append(zstdlow[j]) ; zstdupplst.append(zstdupp[j])         ###
    Appending relevant entries from z lower- and upper boundaries-data
71         k = k + 1
72     print(specnames[i], k)
73     if k == 0:
74         zlst.append(1.0)
    ###Appending relevant entries from z-data

```

```

75     zstdlowlst.append(0.26) ; zstdupplst.append(0.36)           ###Appending relevant
    entries from z lower- and upper boundaries-data
76     else:
77         print("no randomness for you!")
78
79     ###Converting lists to arrays: McKean&Tjoa data.
80     Smaxarr = np.asarray(Smaxlst)
81     vmaxarr = np.asarray(vmaxlst)
82     Smaxerr = np.asarray(Smaxstdlst)
83     vmaxerr = np.asarray(vmaxstdlst)
84     zarr = np.asarray(zlst)
85     zerrlow = np.asarray(zstdlowlst) ; zerrupp = np.asarray(zstdupplst)
86
87
88     #####Synthesis II: Reloaded. Computation of results.
    #####
89     print(">>> Synthesis II: Reloaded")
90
91     B = 50e-6 #Gauss ###Estimate for magnetic field strength of source
92
93     def unredshiftinator(z, vmaxobs):    ###Function to unshift the redshifted peak
    frequencies
94         return (1+z)*vmaxobs
95
96     def angsize(S, z, v):                ###Function to compute angular size in mas
97         S = S/1000
98         return ((8**(5/4))*(B**(1/4))*(S**(1/2))*((1+z)**(1/4)))/(v**(5/4))
99
100    def distance(z):                      ###Function to compute distance in Mpc
101        return np.log(1+z)*299792.458/67.80
102
103    def linsize(theta, D):               ###Function to compute linear size in pc
104        theta = (((theta/1000)/60)/60)*np.pi/180
105        return 1000000*np.tan(theta)*D
106
107    def angerr(S, z, vobs, Serr, zerr, vobserr):    ###Uncertainty function for angular size.
108        C = (8**(5/4))*(B**(1/4))
109        dadS = C*((1+z)**(1/4))/(2*(vobs**(5/4))*(S**(1/2)))
110        dadz = C*(S**(1/2))/(4*(vobs**(5/4))*((1+z)**(3/4)))
111        dadvobs = -C*(5*(S**(1/2))*((1+z)**(1/4)))/(4*(vobs**(9/4)))
112        return np.sqrt((dadS*Serr)**2 + (dadz*zerr)**2 + (dadvobs*vobserr)**2)
113
114    def plotx(S, v):
115        return ((S/1000)**0.5)*(v**-1.25)
116
117    def plotxerr(S, v, Serr, verr):
118        dxdS = (0.5*(1000*S)**-0.5)*(v**-1.25)
119        dxdv = -1.25*((S/1000)**0.5)*(v**-2.25)
120        return np.sqrt((dxdS*Serr)**2 + (dxdv*verr)**2)
121
122    ###Computing relevant data arrays using previously defined functions: McKean&Tjoa data.
123    vmaxemit = unredshiftinator(zarr, vmaxarr)
124    angmar = angsize(Smaxarr, vmaxarr, zarr)
125    angerrlow = angerr(Smaxarr, zarr, vmaxarr, Smaxerr, zerrlow, vmaxerr)
126    for i in range(len(angerrlow)):
127        if angerrlow[i] > angmar[i]:
128            angerrlow[i] = angmar[i]
129    angerrupp = angerr(Smaxarr, zarr, vmaxarr, Smaxerr, zerrupp, vmaxerr)
130    lindist = np.zeros(len(angmar))
131    extdist = np.zeros(len(angmar))
132    for i in range(len(zarr)):
133        entry = cosmocalc(zarr[i])
134        lindist[i] = entry[1]
135        extdist[i] = entry[0]
136    linext = linsize(angmar, extdist)
137
138    ###Printing table of relevant statistics: McKean&Tjoa data.
139    print("\033[1m", "Name", "z", "Lum. Distance", "Smax", "vmax", "vmax emit*ang.",
    size*lin.ext. ", "\033[0m")
140    print("\033[1m", " ", " ", " (Mpc)", " (mJy)", " (GHz)", " (GHz)", " (mas)

```

```

141     * (pc)", "\033[0m")
142 for i in range(len(angmar)):
143     print("{0} {2} {3} {4} {5} {6} {7} {8}".format(names[i], " N/A ", np
144           .round(zarr[i],3), np.round(lindist[i],1), np.round(Smaxarr[i],1), np.round(vmaxarr[i
145           ],2), np.round(vmaxemit[i],2), np.round(angmar[i],2), np.round(angerrlow[i],2), np.round
146           (angerrupp[i],2), np.round(linext[i],1)))
147
148 angmarGPS = []
149 angerrlowGPS = [] ; angerruppGPS = []
150 angmarGPPL = []
151 angerrlowGPPL = [] ; angerruppGPPL = []
152 for i in range(len(typelst)):
153     if fm.fnmatch(typelst[i], "GPS") is True:
154         angmarGPS.append(angmar[i])
155         angerrlowGPS.append(angerrlow[i]) ; angerruppGPS.append(angerrupp[i])
156     elif fm.fnmatch(typelst[i], "GPPL") is True:
157         angmarGPPL.append(angmar[i])
158         angerrlowGPPL.append(angerrlow[i]) ; angerruppGPPL.append(angerrupp[i])
159 angmarGPS = np.asarray(angmarGPS)
160 angerrlowGPS = np.asarray(angerrlowGPS) ; angerruppGPS = np.asarray(angerruppGPS)
161 angmarGPPL = np.asarray(angmarGPPL)
162 angerrlowGPPL = np.asarray(angerrlowGPPL) ; angerruppGPPL = np.asarray(angerruppGPPL)
163
164 ###Creating the data to be plotted on the x-axis.
165 plotdatGPS = plotx(Smax1dat, vmax1dat)
166 plotdatGPPL = plotx(Smax2dat, vmax2dat)
167 plotdatSNElf = plotx(SmaxdatSNElf, vmaxdatSNElf)
168 plotdatSNElb = plotx(SmaxdatSNElb, vmaxdatSNElb)
169 plotdatSNElc = plotx(SmaxdatSNElc, vmaxdatSNElc)
170
171 ###Uncertainties in our data.
172 plotstdGPS = plotxerr(Smax1dat, vmax1dat, Smax1std, vmax1std)
173 plotstdGPPL = plotxerr(Smax2dat, vmax2dat, Smax2std, vmax2std)
174
175 #####Synthesis III: Revolutions. Plotting the results.
176 #####
177 print(">>> Synthesis III: Revolutions")
178
179 ###Plot no.2: Peak flux versus angular size.
180 fig2 = figure(figsize=(40/2.54, 40/2.54))
181 frame2 = fig2.add_subplot(1,1,1, xscale='log', yscale='log')
182 frame2.set_title("Peak flux density vs angular size", fontsize=20)
183 assymerryGPS = [angerrlowGPS, angerruppGPS]
184 assymerryGPPL = [angerrlowGPPL, angerruppGPPL]
185 frame2.errorbar(plotdatGPS, angmarGPS, xerr=plotstdGPS, yerr=assymerryGPS, ecolor="k",
186               elinewidth=1, marker="*", mfc="r", ms=12, ls="", label="McKean very faint GPS sample")
187 frame2.errorbar(plotdatGPPL, angmarGPPL, xerr=plotstdGPPL, yerr=assymerryGPPL, ecolor="k",
188               elinewidth=1, marker="*", mfc="y", ms=12, ls="", label="McKean very faint GPS+PL sample")
189 frame2.errorbar(plotdatSNElf, angmarSNElf, ecolor="k", elinewidth=1, marker="*", mfc="g", ms
190               =12, ls="", label="Snellen's faint GPS sample")
191 frame2.errorbar(plotdatSNElb, angmarSNElb, ecolor="k", elinewidth=1, marker="*", mfc="c", ms
192               =12, ls="", label="Snellen's bright GPS sample")
193 frame2.errorbar(plotdatSNElc, angmarSNElc, ecolor="k", elinewidth=1, marker="*", mfc="b", ms
194               =12, ls="", label="Snellen's CSS sample")
195 frame2.set_xlabel('$S_{max}^{0.5} * v_{max}^{-1.25}$ (Jy$^{0.5}$GHz$^{-1.25}$)', fontsize
196               =16)
197 frame2.set_ylabel('Angular size (mas)', fontsize=16)
198 frame2.tick_params(labelsize=15)
199 #frame2.legend(loc=0)
200
201 show()

```

## B.7 Differential number density calculator

The following script is used to synthesize results to obtain differential number densities.

Scripts/Thesis\_synthesis8\_v6-3.py

```

1 #!/usr/bin/env python
2 from __future__ import division
3 import numpy as np
4 import numpy.random as npr
5 import scipy.optimize as spo
6 import fnmatch as fm
7 from matplotlib.pyplot import figure, show
8 #####THIS SCRIPT COMPUTES DIFFERENTIAL NUMBER COUNTS. SOMETHING IS AMISS!
9
10
11 #####Synthesis I: Reading Data.
12 #####
13 print(">>> Synthesis I")
14
15 #####Reading McKean&Tjoo's name and data (because it's largely monocolumnar, this takes
16 some time and messy coding).
17 dat0TJOA = np.genfromtxt("poslist-McKean.dat", dtype=None, usecols=[0], comments="#")
18 dat1TJOA = np.genfromtxt("poslist-McKean.dat", dtype=None, usecols=[1], comments="#")
19 dat2TJOA = np.genfromtxt("poslist-McKean.dat", dtype=None, usecols=[2], comments="#")
20
21 namesTJOA, RAhrsTJOA, decdegTJOA = np.split(dat0TJOA, 3)
22 namesTJOA = namesTJOA.astype(str).flatten()
23 RAhrsTJOA = np.asarray(RAhrsTJOA.astype(float).flatten())
24 decdegTJOA = np.asarray(decdegTJOA.astype(float).flatten())
25
26 poep1TJOA, RAminTJOA, decminTJOA = np.split(dat1TJOA, 3)
27 RAminTJOA = np.asarray(RAminTJOA.astype(float).flatten())
28 decminTJOA = np.asarray(decminTJOA.astype(float).flatten())
29
30 poep2TJOA, RAssecTJOA, decsecTJOA = np.split(dat2TJOA, 3)
31 RAssecTJOA = np.asarray(RAssecTJOA.astype(float).flatten())
32 decsecTJOA = np.asarray(decsecTJOA.astype(float).flatten())
33
34 #####Reading McKean&Tjoo's synch and synch+powlaw data for spectral peaks.
35 purenamesTJOA = np.genfromtxt("ThesisTable_synchrotronNEW.dat", dtype=str, usecols=[1],
36 comments="#")
37 combinamesTJOA = np.genfromtxt("ThesisTable_synch-powlawNEW.dat", dtype=str, usecols=[1],
38 comments="#")
39 purevmaxTJOA = np.genfromtxt("ThesisTable_synchrotronNEW.dat", dtype=float, usecols=[10],
40 comments="#")
41 combivmaxTJOA = np.genfromtxt("ThesisTable_synch-powlawNEW.dat", dtype=float, usecols
42 =[14], comments="#")
43 vmaxTJOA = np.concatenate((purevmaxTJOA, combivmaxTJOA))
44 pureSmaxTJOA = np.genfromtxt("ThesisTable_synchrotronNEW.dat", dtype=float, usecols=[12],
45 comments="#")
46 combiSmaxTJOA = np.genfromtxt("ThesisTable_synch-powlawNEW.dat", dtype=float, usecols
47 =[16], comments="#")
48 SmaxTJOA = np.concatenate((pureSmaxTJOA, combiSmaxTJOA))
49
50 #####Reading Snellen's name and data.
51 rawnamesSNEL = np.genfromtxt("Snellenfulltab.dat", dtype=str, usecols=[0], comments="#")
52 rawRAhrsSNEL, rawRAminSNEL, rawRAssecSNEL, rawdecdegSNEL, rawdecminSNEL, rawdecsecSNEL = np.
53 genfromtxt("Snellenfulltab.dat", dtype=float, usecols=[1, 2, 3, 4, 5, 6], unpack=True,
54 comments="#")
55 Sgb5, Svla8, Snvss14 = np.genfromtxt("Snellenfulltab.dat", dtype=float, usecols=[9, 13, 16],
56 unpack=True, comments="#")
57
58 #####Selecting only Snellen data which satisfy McKean's criteria.
59 namesSNEL = []
60 RAhrsSNEL = [] ; RAminSNEL = [] ; RAssecSNEL = []
61 decdegSNEL = [] ; decminSNEL = [] ; decsecSNEL = []
62 k = len(rawnamesSNEL)
63 for i in range(len(Svla8)):
64     print("8.4 GHz flux is:", Svla8[i])
65     if Svla8[i] > 16.5:
66         alpha = (np.log10(Sgb5[i]) - np.log10(Snvss14[i])) / (np.log10(5.0) - np.log10(1.4))
67         print("alpha is:", alpha)
68         if 10.0 > alpha > -0.55:
69             k = k - 1

```

```

59     namesSNEL.append(rawnamesSNEL[i])
60     RAhrsSNEL.append(rawRAhrsSNEL[i])
61     RAminSNEL.append(rawRAminSNEL[i])
62     RAsecSNEL.append(rawRAsecSNEL[i])
63     decdegSNEL.append(rawdecdegSNEL[i])
64     decminSNEL.append(rawdecminSNEL[i])
65     decsecSNEL.append(rawdecsecSNEL[i])
66     elif alpha == np.inf:
67         print(rawnamesSNEL[i], "missed either its 5 or 1.4 GHz entry.")
68     else:
69         print(rawnamesSNEL[i], "had a spectral index between 5 and 1.4 GHz steeper
70         than -0.5.")
71     else:
72         print(rawnamesSNEL[i], "had a 8.6 GHz flux which was too low to register.")
73 print("Number of rejected entries:", k)
74 RAhrsSNEL = np.asarray(RAhrsSNEL) ; RAminSNEL = np.asarray(RAminSNEL) ; RAsecSNEL = np.
75     asarray(RAsecSNEL)
76 decdegSNEL = np.asarray(decdegSNEL) ; decminSNEL = np.asarray(decminSNEL) ; decsecSNEL =
77     np.asarray(decsecSNEL)
78
79 #####Reading Snellen's synch data for spectral peaks.
80 peaknamesSNEL = np.genfromtxt("Snellenpeaktab.dat", dtype=str, usecols=[0], comments="#")
81 vmaxSNEL = np.genfromtxt("Snellenpeaktab.dat", dtype=float, usecols=[1], comments="#")
82 SmaxSNELraw = np.genfromtxt("Snellenpeaktab.dat", dtype=float, usecols=[2], comments="#")
83
84 #####Synthesis II: Reloaded. Computation of results.
85 #####
86 print(">>> Synthesis II: Reloaded")
87
88 def posmaker(RAhrs, RAmin, RAsec, decdeg, decmin, decsec): ###Function that computes the RA
89     and dec positions in fractional hours and degrees respectively for plotting purposes
90     RApos = RAhrs + RAmin/60 + (RAsec/60)/60
91     decpos = decdeg + decmin/60 + (decsec/60)/60
92     return RApos, decpos
93
94 def areacalc(RAlow, RAupp, declow, decupp):
95     declow = declow*np.pi/180 ; decupp = decupp*np.pi/180
96     RAlow = 15*RAlow*np.pi/180 ; RAupp = 15*RAupp*np.pi/180
97     return (RAupp - RAlow)*(np.sin(decupp) - np.sin(declow))
98
99 def difnumdensflux(Slow, Supp, Sp):
100     k = 0
101     reject = 0
102     for i in range(len(Sp)):
103         if Sp[i] >= Slow and Sp[i] <= Supp:
104             k = k + 1
105     return k
106
107 #####Computing the positions of Tjoa&McKean's sources.
108 posTJOA = posmaker(RAhrsTJOA, RAminTJOA, RAsecTJOA, decdegTJOA, decminTJOA, decsecTJOA)
109
110 #####Computing the positions of Snellen's sources.
111 posSNEL = posmaker(RAhrsSNEL, RAminSNEL, RAsecSNEL, decdegSNEL, decminSNEL, decsecSNEL)
112
113 #####Computing positions of GPS samples, McKean&Tjoa.
114 pureRATJOA = []
115 puredecTJOA = []
116 for i in range(len(purenamesTJOA)):
117     for j in range(len(namesTJOA)):
118         if fm.fnmatch(purenamesTJOA[i], "{}".format(namesTJOA[j])) is True:
119             pureRATJOA.append(posTJOA[0][j])
120             puredecTJOA.append(posTJOA[1][j])
121 pureRATJOA = np.asarray(pureRATJOA)
122 puredecTJOA = np.asarray(puredecTJOA)
123 peaknamesTJOA = list(purenamesTJOA)
124
125 combiRATJOA = []
126 combidecTJOA = []
127 for i in range(len(combinamesTJOA)):

```

```

124     for j in range(len(namesTJOA)):
125         if fm.fnmatch(combinamesTJOA[i], "{}".format(namesTJOA[j])) is True:
126             combiRATJOA.append(posTJOA[0][j])
127             combidecTJOA.append(posTJOA[1][j])
128 combiRATJOA = np.asarray(combiRATJOA)
129 combidecTJOA = np.asarray(combidecTJOA)
130 peaknamesTJOA.extend(combinamesTJOA)
131
132 SmaxTJOA = np.concatenate((pureSmaxTJOA,combiSmaxTJOA))
133 peakRATJOA = np.concatenate((pureRATJOA,combiRATJOA))
134 peakdecTJOA = np.concatenate((puredecTJOA,combidecTJOA))
135
136 ###Computing positions of GPS samples, Snellen.
137 SmaxSNEL = []
138 peakRASNEL = []
139 peakdecSNEL = []
140 for i in range(len(peaknamesSNEL)):
141     for j in range(len(namesSNEL)):
142         if fm.fnmatch(peaknamesSNEL[i], "{}".format(namesSNEL[j])) is True:
143             SmaxSNEL.append(SmaxSNELraw[i])
144             peakRASNEL.append(posSNEL[0][j])
145             peakdecSNEL.append(posSNEL[1][j])
146 SmaxSNEL = np.asarray(SmaxSNEL)
147 peakRASNEL = np.asarray(peakRASNEL)
148 peakdecSNEL = np.asarray(peakdecSNEL)
149
150 ###Loop that splits Tjoa&McKean's data (only peaked) into ists two component fields of
151     view and computing areas.
152 RAlowhighTJOA = 17 ; RAupphighTJOA = 18 ; declowhighTJOA = 55 ; decupphighTJOA = 60
153 RAlowlowTJOA = 1 ; RAupplowTJOA = 2 ; declowlowTJOA = 35 ; decupplowTJOA = 45
154 RAhighpTJOA = [] ; dechighpTJOA = [] ; SmaxhighTJOA = []
155 RAlowpTJOA = [] ; declowpTJOA = [] ; SmaxlowTJOA = []
156 for i in range(len(peakRATJOA)):
157     if peakRATJOA[i] > 2.0:
158         RAhighpTJOA.append(peakRATJOA[i])
159         dechighpTJOA.append(peakdecTJOA[i])
160         SmaxhighTJOA.append(SmaxTJOA[i])
161     elif peakRATJOA[i] < 2.0:
162         RAlowpTJOA.append(peakRATJOA[i])
163         declowpTJOA.append(peakdecTJOA[i])
164         SmaxlowTJOA.append(SmaxTJOA[i])
165
166 areahighTJOA = areacalc(RAlowhighTJOA, RAupphighTJOA, declowhighTJOA, decupphighTJOA)
167 arealowTJOA = areacalc(RAlowlowTJOA, RAupplowTJOA, declowlowTJOA, decupplowTJOA)
168 print("Our high field:", areahighTJOA, "sr")
169 print("Our low field:", arealowTJOA, "sr")
170 print("Our full field:", areahighTJOA+arealowTJOA, "sr")
171
172 ###Loop that splits Snellen's data (only peaked) into its two component fields of view
173     and computing areas.
174 RAlowhighSNEL = 15 ; RAupphighSNEL = 20 ; declowhighSNEL = 58 ; decupphighSNEL = 75
175 RAlowlowSNEL = 4 ; RAupplowSNEL = 8.5 ; declowlowSNEL = 58 ; decupplowSNEL = 75
176 RAhighpSNEL = [] ; dechighpSNEL = [] ; SmaxhighSNEL = []
177 RAlowpSNEL = [] ; declowpSNEL = [] ; SmaxlowSNEL = []
178 for i in range(len(peakRASNEL)):
179     if peakRASNEL[i] > 9.0:
180         RAhighpSNEL.append(peakRASNEL[i])
181         dechighpSNEL.append(peakdecSNEL[i])
182         SmaxhighSNEL.append(SmaxSNEL[i])
183     elif peakRASNEL[i] < 9.0:
184         RAlowpSNEL.append(peakRASNEL[i])
185         declowpSNEL.append(peakdecSNEL[i])
186         SmaxlowSNEL.append(SmaxSNEL[i])
187
188 areahighSNEL = areacalc(RAlowhighSNEL, RAupphighSNEL, declowhighSNEL, decupphighSNEL)
189 arealowSNEL = areacalc(RAlowlowSNEL, RAupplowSNEL, declowlowSNEL, decupplowSNEL)
190 print("Snellen's high field:", areahighSNEL, "sr")
191 print("Snellen's low field:", arealowSNEL, "sr")

```

```

192 print("Snellen's full field:", areahighSNEL+arealowSNEL, "sr")
193
194 Smed = np.asarray([30,60,120,240,480])
195 ###Computing number density-peak flux convolution per steradian as a function of peak
    flux (differential method).
196 def peakconvo(Smaxhigh, Smaxlow, areatot):
197     Sbinsize = Smed/3          ###Binsize for number counts
198     Slow = (2/3)*Smed         ###Lower bound of bin
199     Supp = (4/3)*Smed         ###Upper bound of bin
200
201     nsrdifhigh = []
202     nsrdiflow = []
203     for i in range(len(Smed)):    ###Loop that uses function to differentially count
        number of sources per bin
204         nsrdifhigh.append(difnumdensflux(Slow[i], Supp[i], Smaxhigh))
205         nsrdiflow.append(difnumdensflux(Slow[i], Supp[i], Smaxlow))
206     nsrdifhigh = np.asarray(nsrdifhigh)    ###Converting counts in high field to array
207     nsrdiflow = np.asarray(nsrdiflow)      ###Converting counts in low field to array
208
209     nsrdiftot = nsrdifhigh+nsrdiflow        ###Calculating total amount of sources per bin
210
211     densdiftot = nsrdiftot/areatot          ###Computing source density per bin in total
212
213     dndS = nsrdiftot/((Supp-Slow)/1000)    ###dn/dS -> density over binwidth
214     plotypeak = (dndS/areatot)*((Smed/1000)**(5/2))    ###Plotted values on y-axis
215     plotyerr = (np.sqrt(nsrdiftot)/nsrdiftot)*plotypeak    ###Poissonian uncertainties
        in y-axis values
216     ### 0      1      2      3      4
217     return Sbinsize, nsrdiftot, densdiftot, plotypeak, plotyerr
218
219 ###McKean&Tjoa data.
220 TJOA = peakconvo(SmaxhighTJOA, SmaxlowTJOA, areahighTJOA+arealowTJOA)
221 SbinsizeTJOA = TJOA[0]
222 plotypeakTJOA = TJOA[3]
223 plotyerrTJOA = TJOA[4]
224 plotypeak2TJOA = plotypeakTJOA/(Smed**(5/2))
225 plotyerr2TJOA = plotyerrTJOA/(Smed**(5/2))
226 print(TJOA[1])
227
228 ###Snellen data.
229 SNEL = peakconvo(SmaxhighSNEL, SmaxlowSNEL, areahighSNEL+arealowSNEL)
230 SbinsizeSNEL = SNEL[0]
231 plotypeakSNEL = SNEL[3]
232 plotyerrSNEL = SNEL[4]
233 plotypeak2SNEL = plotypeakSNEL/(Smed**(5/2))
234 plotyerr2SNEL = plotyerrSNEL/(Smed**(5/2))
235 print(SNEL[1])
236
237 ###Combined data
238 SbinsizeCOMB = SbinsizeTJOA
239 nsrCOMB = TJOA[1]+SNEL[1]
240 plotypeakCOMB = plotypeakSNEL + plotypeakTJOA
241 plotyerrCOMB = (np.sqrt(nsrCOMB)/nsrCOMB)*plotypeakCOMB
242 plotypeak2COMB = plotypeak2SNEL + plotypeak2TJOA
243 plotyerr2COMB = (np.sqrt(nsrCOMB)/nsrCOMB)*plotypeak2COMB
244 print(nsrCOMB)
245
246 ###De Vries data.
247 SmedVRIES = 4000
248 SbinsizeVRIES = 2000.0
249 plotypeakVRIES = 22.0
250 plotyerrVRIES = 10.0
251 plotypeak2VRIES = plotypeakVRIES/(SmedVRIES**(5/2))
252 plotyerr2VRIES = plotyerrVRIES/(SmedVRIES**(5/2))
253
254 with open("ThesisTable_numdifdensTJOA2.dat", "a") as myfile:
255     for i in range(len(nsrCOMB)):
256         myfile.write("{0} & {1} & {2} & {5}$\pm${6} & {11}$\pm${12} \\\\".format(Smed[i]
        ], Smed[i]/3, TJOA[1][i], SNEL[1][i], nsrCOMB[i], np.round(plotypeakTJOA[i], 3), np.round(
        plotyerrTJOA[i], 3), np.round(plotypeakSNEL[i], 3), np.round(plotyerrSNEL[i], 3), np.round

```

```

        (plotypeakCOMB[i],3),np.round(plotyerrCOMB[i],3),np.round(plotypeak2TJOA[i],10),np.
        round(plotyerr2TJOA[i],10),np.round(plotypeak2SNEL[i],10),np.round(plotyerr2SNEL[i
        ],10),np.round(plotypeak2COMB[i],10),np.round(plotyerr2COMB[i],10)))
257         myfile.write("\n")
258
259
260 #####Synthesis III: Revolutions. Plotting the results.
261 #####
262 print(">>> Synthesis III: Revolutions")
263
264 ###Plot no.0: number density/peak convolution as a function of peak flux.
265 fig0 = figure(figsize=(40/2.54, 40/2.54))
266 frame0 = fig0.add_subplot(1,1,1,xscale='log',yscale='log')
267 frame0.errorbar(Smed,plotypeakTJOA,xerr=SbinsizeTJOA,yerr=plotyerrTJOA,ecolor="k",
268                 elinewidth=2,marker="*",mfc="r",ms=30,ls="",label="McKean's very faint GPS sample")
269 frame0.errorbar(Smed,plotypeakSNEL,xerr=SbinsizeSNEL,yerr=plotyerrSNEL,ecolor="k",
270                 elinewidth=2,marker="*",mfc="y",ms=30,ls="",label="Snellen's GPS sample")
271 #frame0.errorbar(Smed,plotypeakCOMB,xerr=SbinsizeCOMB,yerr=plotyerrCOMB,ecolor="k",
272                 elinewidth=2,marker="*",mfc="g",ms=20,ls="",label="Total GPS sample")
273 frame0.errorbar(SmedVRIES,plotypeakVRIES,xerr=SbinsizeVRIES,yerr=plotyerrVRIES,ecolor="k",
274                 elinewidth=2,marker="*",mfc="c",ms=20,ls="",label="De Vries' sample")
275 frame0.set_xlabel("$S_{max}$ (mJy)",fontsize=35)
276 frame0.set_ylabel("$dN(>S)/dS*S^{5/2}$ (Sr^{-1}Jy^{3/2})",fontsize=35)
277 frame0.tick_params(labelsize=30)
278
279 ###Plot no.1: number density as a function of peak flux.
280 fig1 = figure(figsize=(40/2.54, 40/2.54))
281 frame1 = fig1.add_subplot(1,1,1,xscale='log',yscale='log')
282 frame1.errorbar(Smed,plotypeak2TJOA,xerr=SbinsizeTJOA,yerr=plotyerr2TJOA,ecolor="k",
283                 elinewidth=2,marker="*",mfc="r",ms=30,ls="",label="McKean's very faint GPS sample")
284 frame1.errorbar(Smed,plotypeak2SNEL,xerr=SbinsizeSNEL,yerr=plotyerr2SNEL,ecolor="k",
285                 elinewidth=2,marker="*",mfc="y",ms=30,ls="",label="Snellen's GPS sample")
286 #frame1.errorbar(Smed,plotypeak2COMB,xerr=SbinsizeCOMB,yerr=plotyerr2COMB,ecolor="k",
287                 elinewidth=2,marker="*",mfc="g",ms=20,ls="",label="Total GPS sample")
288 frame1.errorbar(SmedVRIES,plotypeak2VRIES,xerr=SbinsizeVRIES,yerr=plotyerr2VRIES,ecolor="k",
289                 elinewidth=2,marker="*",mfc="c",ms=20,ls="",label="De Vries' sample")
290 frame1.set_xlabel("$S_{max}$ (mJy)",fontsize=35)
291 frame1.set_ylabel("$dN(>S)/dS$ (Sr^{-1}Jy^{-1})",fontsize=35)
292 frame1.tick_params(labelsize=30)
293
294 show()

```

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