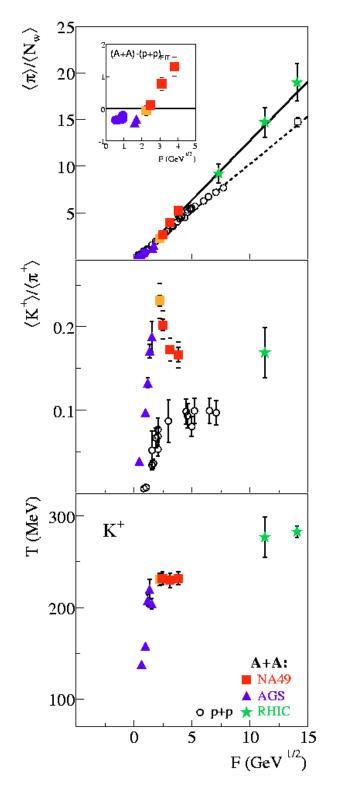
Energy dependence of transverse mass spectra of kaons produced in p+p and p+p interactions <u>A compilation</u>

Benjamin Lungwitz and Michael Kliemant

- Anomaly observed in the T vs $\sqrt{s_{NN}}$ dependence for central Pb+Pb collisions
- How does this dependence look like for p+p interactions?

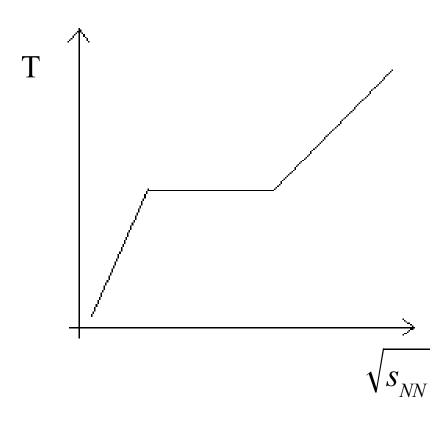
See also in our paper: arXiv:hep-ex/0308002



 Pion multiplicity a compilation of p+p data existed

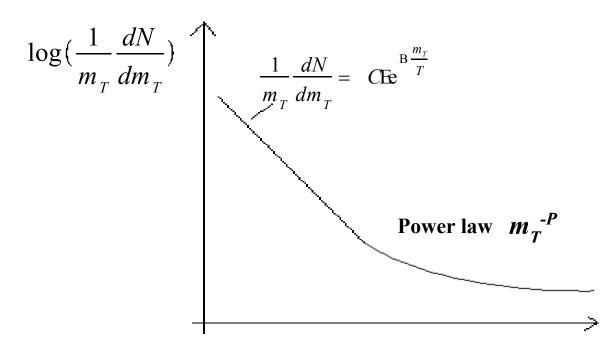
• Kaon multiplicity a compilation of p+p data existed

 Kaon inverse slope parameter a compilation of p+p data did not exist

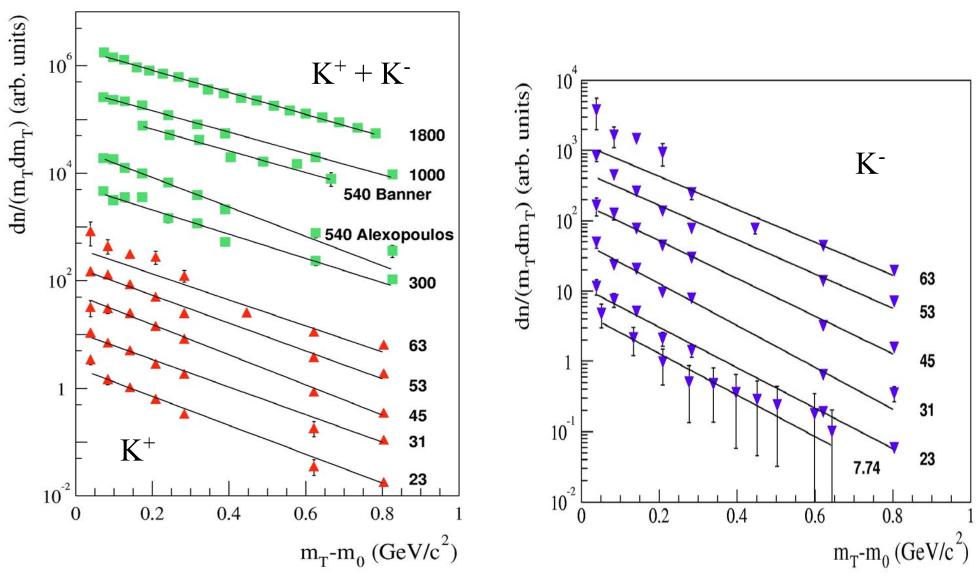


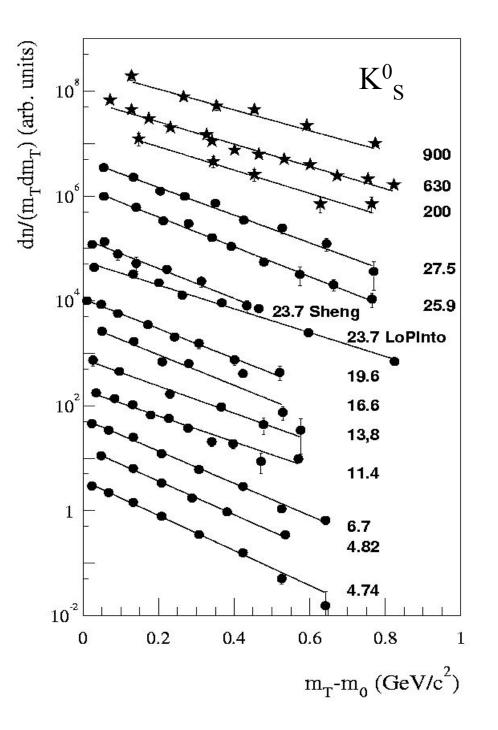
- The Pb+Pb results show
 a phase transition like
 behaviour (□caloric curve
 □Kochtopfmodel□)
- Do we see a similar dependence for p+p?

- We compiled and analyzed data on m_T spectra of K^0_S , K^+ and K^- in p+p, p+ \overline{p} interactions at all energies ($\sqrt{s_{NN}} = 4.74 \text{GeV} 1.8 \text{TeV}$). The data originate fom fixed target and collider experiments performed during the last 30 years.
- The spectra were fitted by a simple exponential function
- This parametrization is valid only in low m_T region ($m_T < 1.2 \text{GeV}$); at higher m_T the power law behaviour m_T^{-P}



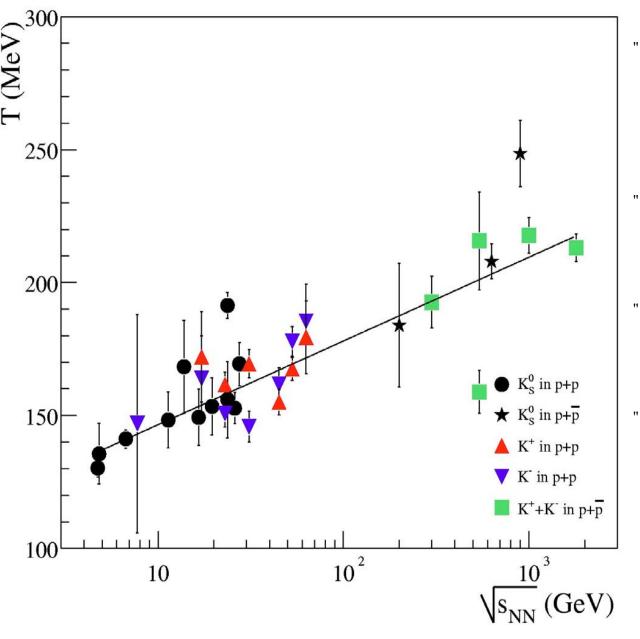
$dn/(m_T dm_T)$ spectra





- Exponential function fits the data
- Inverse slope parameter
 increases with energy

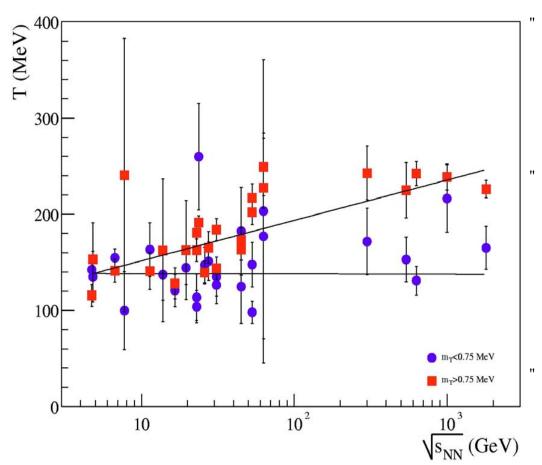
Fits in full $m_T (m_T < 1.2 \text{ GeV})$ region



- No obvious difference between K⁰_S, K⁺ and K⁻
- T for p+p and $p+\overline{p}$ reactions is similar
- Logarithmic increase of T with $\sqrt{s_{NN}}$
- Few points do not match the fit (possible systematic errors)

Fits in low and high m_T regions

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Is the T parameter the same when we fit different m_T regions ?

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In the low m_T region ($m_T < 0.75$) T seems to be energy independent

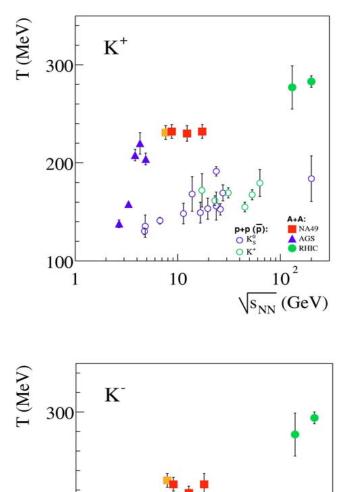
In the high m_T region (m_T >0.75 GeV/c²) it shows logarithmic increase (beginning of the power law ?)

Possible large systematic errors when using different m_T intervals for a comparison

For a comparison with A+A data a similar m_T range was used

Comparison p+p to Pb+Pb

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10

s_{NN} (GeV)

200

100

1

- No significant difference between
 K⁺ and K⁻ dependences for both p+p
 and Pb+Pb collisions
 - T(p+p) < T(Pb+Pb) at all energies
- No significant \Box transition like structure in T(p+p)
- Quality of the world p+p data is
 poor → new measurements are
 needed

