



Goals:

- Estimation of dE/dx separation powers.
- dE/dx Calibration elements and dE/dx ID efficiency/fake rate measurements.
- Investigation of different gas mixtures using FastSim & Garfield simulation.
- Estimation of IP resolution.

Using the FastSim 0.1.3 version we simulated different decays of B/Bbar to obtain a pure samples of analyzed particles covering as much as possible wide momentum range:

- **Electrons and Muons from J/ψ decays;**
- Pions from B decays;
- > Protons from B and J/ ψ decays.

For dE/dx analysis we selected samples of electrons, muons, pions, kaons and protons without particles identification criteria.

Pions Selection



Decay: $B^0 \rightarrow \pi^+\pi^-$, $B^0 bar \rightarrow \pi^+\pi^-\pi^+\pi^-$

TRKDocaXY_xy < 0.04 m |TRKDocaXY_z| < 0.2 m



Kaons Selection



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Simulation of events with "only" Kaons in the final state.



Selected Samples For dE/dx Analysis





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Separation Powers of Electrons and Other Particles



SuperB

Separation Powers of Muons And Other Particles



SuperB

Pion / Kaon And Pion / Proton Sepatation Powers.



Kaon / Proton Separation Power



Particles Separation for SuperB DCH

dE/dx Tuning



dE/dx Tuning is very important for data, but should be tested on MC info.



We parameterized dE/dx as a function of $\beta\gamma$ for muons, pions, kaons and protons simultaneously.

We fit three $\beta\gamma$ regions ([0. - 1.0], [1.0 - 4.5] and [4.5 - 50.]) to a power law + third order polynomial.

$$dE / dx (\beta \gamma) = A \cdot (\beta \gamma)^n + B_0 + B_1 \cdot (\beta \gamma) + B_2 \cdot (\beta \gamma)^2 + B_3 \cdot (\beta \gamma)^2$$

At very large βγ dE/dx is constant ("Fermi Plateau")

In the region $\beta\gamma > 50$ (electrons) for dE/dx parameterization exp-function can be used.

$$dE / dx(\beta \gamma) = C \cdot e^{D \cdot (\beta \gamma)} + E$$

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dE/dx Tuning, cntd.



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IE/dx Tuning, cntd.



g dE/dx

For particle identification using dE/dx the variables σ was constructed. The value of σ is given by:

$$\sigma = \frac{(dE / dx)_{Measured} - (dE / dx)_{Expected}}{\sigma_{dE / dx}}$$

dE/dx is a function of $\beta\gamma$ for each particle. For each type of particles (e, m, p, K, p) using $\beta\gamma$ we find the expected value of (dE/dx)_{Expected}.

The track level quantity dE/dx can be thought of as the total charge (Q) deposited by the track divided by the total path length of the track.

The error on Q should be proportional to $Q^{1/2}$ and therefore the resolution of $\sigma_{dE/dx}$ should be given by equation:

$$\sigma_{dE/dx} \approx \frac{Q^{1/2}}{L} = \frac{((dE/dx) \cdot L)^{1/2}}{L} = \frac{(dE/dx)_{Measured}^{1/2}}{L^{1/2}} L = N_{hit}/\sin(\theta) \text{ is a path length.}$$
We plan to parametrize the dE/dx resolution
in terms of three functions:
$$\sigma_{dE/dx} = \frac{f[(dE/dx)_{Measured}^{1/2}]}{g[N_{Hit}^{1/2}] \cdot h[1/\sin^{1/2}(\theta)]}$$



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dE/dx Identification, cntd.





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dE/dx Identification Efficiency Measurement



KID efficiency: $\varepsilon = \#Kaons(|\sigma_K|<3) / \#Kaons(Before ID Cut)$ π ID efficiency: $\varepsilon = \#Pions(|\sigma_{\pi}|<3) / \#Pions(Before ID Cut)$ Fake Rate (K fake as π): f.r. = $\#Kaons(|\sigma_{\pi}|<3) / \#Kaons(Before ID Cut)$ Fake Rate (π fake as K): f.r. = $\#Pions(|\sigma_K|<3) / \#Pions(Before ID Cut)$



DCH Materials List in FastSim

Currently the following numbers of DCH materials are in the MaterialsList.data

| | Density | Z _{eff} | A _{eff} | | X0 | λ _I |
|------------|------------|-------------------------|------------------|---|-----------|----------------|
| dch-He-lbu | 6.408E- 04 | 23.8 | 46.1 | 0 | 51.16 | 75.65 |
| dch-Wires | 6.237E+00 | 29.0 | 62.4 | 0 | 15.31 | 118.56 |



$$C_4H_{10}$$
: $A_{Iso} = 4 A_C + 10 A_H$
 C_3H_8 : $A_{Pro} = 3 A_C + 8 A_H$

 $Z_{Iso} = 4 Z_{C} + 10 Z_{H}$ $Z_{Pro} = 3 Z_{C} + 8 Z_{H}$

Atomic and Nuclear Properties of Materials http://pdg.lbl.gov/2009/AtomicNuclearProperties/

| Material | Density g/cm ³ | Atomic Number, Z | Atomic Mass, A | Z/A | Rad.Len. X0, g/cm ² | Nucl.Int.L. λ_{i} , g/cm ² |
|-----------|------------------------------|---------------------|-------------------|----------|-----------------------------------|---|
| Н | - | 1 | 1.00794 | - | - | - |
| С | - | 6 | 12.0107 | - | - | - |
| Не | 1.66E-04 | 2 | 4.0026 | 0.499675 | 94.32 | 71.0 |
| Isobutane | 2.49E-03 | 34 | 58.1222 | 0.584974 | 45.23 | 77.1 |
| Propane | 1.87E-03 | 26 | 44.0956 | 0.589628 | 45.37 | 76.7 |

Calculation of the MaterialsList.data DCH Components



Effective Z/A ratio for gas mixtures were calculated as:

$$\left(\frac{Z}{A}\right)_{Eff} = \sum_{i} \omega_{i} \cdot \frac{Z_{i}}{A_{i}} = \sum_{i} \frac{f_{i} \cdot \rho_{i}}{\rho} \cdot \frac{Z_{i}}{A_{i}} \qquad \rho = \sum_{i} f_{i} \cdot \rho_{i}$$

 f_i – is a fraction of gas components, ρ – is the density of the mixture. Radiation and Nuclear interaction lengths for gas mixtures were calculated as:

$$\frac{1}{X_{0}} = \sum_{i} \frac{f_{i} \cdot \rho_{i}}{\rho} \cdot \frac{1}{X_{0i}} \qquad \frac{1}{L} = \sum_{i} \frac{f_{i} \cdot \rho_{i}}{\rho} \cdot \frac{1}{L_{i}}$$

Atomic masses and Atomic Numbers for a gas mixtures were computed as:

$$A_{He / Iso} = \frac{f_{He} \cdot \rho_{He} \cdot A_{He} + f_{Iso} \cdot \rho_{Iso} \cdot A_{Iso}}{\rho_{He / Iso}}$$

 $Z_{He / Iso} = A_{He / Iso} \cdot \left(\frac{Z}{A}\right)_{Eff}$

| Gas Mixture | Density | Atomic | Atomic | Rad.Len. | Nucl.Int.L. |
|---------------------|-------------------|-----------|---------|------------------------------|-------------------------|
| | g/cm ³ | Number, Z | Mass, A | X0, g/cm ² | $\lambda_{l}, g/cm^{2}$ |
| He(60%) & Prop(40%) | 8.476E-04 | 22.81 | 39.38 | 48.32 | 75.98 |
| He(70%) & Iso(30%) | 8.632E-04 | 29.16 | 50.84 | 48.64 | 76.22 |
| He(80%) & Iso(20%) | 6.308E-04 | 26.50 | 46.73 | 50.80 | 75.73 |
| He(90%) & Iso(10%) | 3.984E-04 | 20.92 | 37.83 | 56.20 | 74.69 |

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Comparison of Separation Powers for Different Gas Mixtures



Comparison of Separation Powers for Different Gas Mixtures





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Requirements to the dN/dx Technique



Pulses from electrons belonging to different clusters should not overlapping in time.

A low pulse density in the time gate (prevents possible overlaps).

Nclusters vs particle energy. Gas: He / C4H10



| He/ C ₄ H ₁₀ , % | <n> Clusters</n> | σ |
|--|------------------|-----------------------------------|
| 100 / 0 | 4.28 ± 0.08 | 1.48 ± 0.05 |
| 95 / 05 | 8.74 ± 0.07 | $\textbf{2.68} \pm \textbf{0.05}$ |
| 90 /10 | 13.31 ± 0.07 | $\textbf{3.36} \pm \textbf{0.05}$ |
| 85 /15 | 18.05 ± 0.08 | 4.10 ± 0.06 |
| 80 /20 | 22.56 ± 0.10 | $\textbf{4.49} \pm \textbf{0.07}$ |
| 75 /25 | 27.35 ± 0.11 | 5.01 ± 0.08 |
| 70 /30 | 31.68 ± 0.10 | $\textbf{5.52} \pm \textbf{0.07}$ |

 He/C_4H_{10} (90%/10%) is preferable for the **Cluster Counting technique**

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Comparison of dE/dx & dN/dx Separation Powers



Separation power for ideal cluster counting increased in respect with dE/dx method from ~1.4 σ to 3.2 σ (in the momentum region 2 GeV/c < P < 3 GeV/c).



Requirements to the dE/dx Technique

Want a faster gas mixture with more primary clusters.



As Lorentz angle increases, electron drift path curvature also increases. Want to minimize Lorentz angle for straighter path and larger correlation between distance and time.



Impact Parameter Resolution Estimation (Very Preliminary)



Drift Tube



Cluster counting can contribute to reduce the impact parameters resolution for small impact parameters.

Using Garfield we simulated tracks of minimum ionizing pions crossing drift tube with the impact parameters: b=0., 0.05, 0.1, 0.15, 0.2, 0.25, 0.30, 0.35, 0.40, 0.45 cm.

For each impact parameter we simulated $\sim 2x10^4$ tracks.

mpact Parameter Resolution Estimation Method





The impact parameter represents a systematic overestimate of b by the quantity Δb , usually referred to as the ionization statistics contribution to the impact parameter resolution:

$$\Delta \mathbf{b} = \mathbf{d} - \mathbf{b}$$

We used clusters with $\lambda/4 > \lambda_1 > 0$. cm

Impact Parameter Resolution (Very Preliminary)

Looks like that Cluster Counting method can improve spatial resolution by factor 2-3 in respect with CLEO/BaBar resolutions.

But spatial resolution needs more analysis. In particular:

According to the 4Lol, most of the improvement in impact parameter resolution is obtained by using only the first two clusters, with smaller and smaller contributions from the successive ones.

In this analysis we took into account all ionization clusters.

We need more motivated selection of clusters for estimation of the impact parameter resolution.

 This work is in progress.

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Summary:

- FastSim & Garfield simulation study for different gas compositions was performed.
- Particles separation powers based on FastSim were estimated.
- We performed MaterialsList calculations for different gas compositions for FastSim to estimate particles separation powers for different gas mixtures.
- We tested possible dE/dx calibration procedure in Fast Simulation: parameterized dE/dx as a function of $\beta\gamma$ for muons, pions, kaons and protons simultaneously.
- Based on FastSim we estimated dE/dx Identification efficiencies and Fake Rates for pions and kaons in wide momentum region.
- Using Garfield we estimated different gas characteristics: Drift time, velocity, Lorentz angle, etc. to compare the requirements of dE/dx and dN/dx techniques.
- Preliminary estimation of spatial resolution was presented.
- He/Ibu mixture with He fraction more that 80% and less than 90% can be considered as a candidate of SuperB DCH gas acceptable for both, dE/dx and dN/dx technique (with small size inner cells).