## Dual-Energy Decomposition Methods for Accurate Material Discrimination

ADSA 16

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## ( $\rho_{\mathrm{e}}, \mathrm{Z}_{\mathrm{e}}$ ) is useful for precise and accurate material discrimination

Physical properties provide a system independent feature space.

Different dual-energy decomposition methods available.


Propose an adaptive method for a system-to-system transfer function.

Reduced artifacts and errors over entire material range may lead to increased $P_{D}$ and lowered $P_{F A}$.

Absolute relative error of LAC estimate at 44 keV

| Basis <br> Type | HDPE <br> $\left(Z_{e}=5.4\right)$ | Water <br> $\left(Z_{e}=7.45\right)$ | $A l$ <br> $\left(Z_{e}=13\right)$ | Brominated Delrin <br> $\left(Z_{e}=17.0\right)$ |
| :--- | :--- | :--- | :--- | :--- |
| LAC (PCD) | $0.10 \%$ | $0.18 \%$ | $0.03 \%$ | $3.28 \%$ |
| Adaptive | $0.16 \%$ | $0.14 \%$ | $0.10 \%$ | $0.17 \%$ |

Dual-Energy:
Reduces artifacts


## Approximate Multi-Spectral CT Model

Multi-spectral decomposition approximates energy- and spatial- dependent linear attenuation coefficients (LAC) with a separable basis expansion:



Effectiveness of multi-spectral decomposition depends on the accuracy of the above basis expansion

## Pros and Cons of Two Common Bases

Compton/ photoelectric basis

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Alvarez and Macovski, Phys. Med. Biol., 1976
Ying, Naidu, \& Crawford, J. X-ray Sci. and Tech., 2006
Azevedo et al., IEEE Trans. Nucl. Sci., 2016
```

Pros

1. Works well for materials with effective-Z between 5 and 18
2. Relative absolute error is $<1.3 \%$
3. Mostly commonly used and well-understood

Cons

1. Does not correctly measure the Compton or photoelectric coefficients
2. Photoelectric coefficient is noisy, which limits it practical use

## Basis Materials

Pros

1. Coefficients have some physical relevance, e.g., C and Al
2. Interpolates well within narrow range, relative error is $<0.4 \%$
3. Noise levels of coefficients roughly match noise levels of measured data

Cons

1. Error increases rapidly outside narrow range

## Principal Component Analysis (PCA) Basis

- Effectiveness of multi-energy basis approximation driven by the following approximation

$$
\mu(E, x) \approx \sum_{i} b_{i}(E) c_{i}(x)
$$

- We may choose a basis by minimizing the error in the above approximation for a specified range of materials
- Principal Component Analysis enables us to choose the best basis possible because it minimizes the Mean Square Error
- 3-element PCA basis derived from LAC values of $Z=6$, $7,8,13,26,36$; relative absolute error is $<1 \%$ from $Z$ range 5 to 35
- While Compton, photoelectric, iron bases error is < 3\% from $Z$ range 5 to 35


## Transformation to LAC/ Mono-Energy Basis

LAC basis is simply a linear transformation of any given basis, e.g., Compton/ photoelectric, material basis, etc

To use LAC basis, one must choose a basis and a pair of energies, $\mathrm{E}_{1}$ and $\mathrm{E}_{2}$

Then basis coefficients are LAC values at $\mathrm{E}_{1}$ and $\mathrm{E}_{2}$

$$
\begin{gathered}
\mu(E, x) \approx b_{1}(E) \mu_{1}(x)+b_{2}(E) \mu_{2}(x) \\
\mu(54, x) \approx 0 \mu_{1}(x)+1 \mu_{2}(x) \\
\mu(95, x) \approx 1 \mu_{1}(x)+0 \mu_{2}(x)
\end{gathered}
$$

No matter what basis is used, transformation to LAC basis gives you:

- Basis functions that are unitless
- Basis coefficients that are in LAC units $\left(\mathrm{mm}^{-1}\right)$


## Comparison of Compton/ photoelectric and LAC basis results with Real Measured Data



## Adaptive Dual-Energy Decomposition

Form multiple LAC bases from different basis representations, e.g. bases composed of all pairs of elements that could be present in luggage

For each dual-energy measurement, perform multi-energy decomposition with each LAC basis and choose the result with the lowest residual error

Thus each ray-sum measurement pair could be decomposed using a different set of basis materials

LAC basis is necessary because it maps all bases to a common unit system that can be reconstructed; you cannot reconstruct if each measurement pair lies in a different space

## Dual-Energy Decomposition Experiment with Simulated Polychromatic Data

Spectrum 1:
100 kV , 2 mm Al filter
Spectrum 2:
$160 \mathrm{kV}, 2 \mathrm{~mm} \mathrm{Al}, 2 \mathrm{~mm}$ Cu filters

all objects have 1" diameter

## Dual Energy Decomposition Results

Absolute relative error of LAC estimate at 44 keV

| Basis Type | HDPE <br> $\left(Z_{e}=5.4\right)$ | Water <br> $\left(Z_{e}=7.45\right)$ | AI $\left(Z_{e}=13\right)$ | Brominated <br> Delrin <br> $\left(Z_{e}=17.0\right)$ |
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LAC basis derived from PCA performed on par with
LAC basis derived from Compton/ photoelectric basis

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Dual-Energy:
Reduces artifacts


## Questions?

## Backup Slides

## Multi-Spectral CT for Baggage Inspection



Dual Energy
Reconstruction

## Current methods

- work well for materials with a narrow range of atomic number
- do not provide results with specified units

We develop new multi-spectral methods that produce results with specified units and are accurate for materials with a wider range of atomic number


## Common Dual-Energy Models

## Compton/ Photoelectric Basis

Compton basis function is based on the total Klein-Nishina cross section and photoelectric basis function is $\mathrm{E}^{-3}$

Basis Materials
Basis functions are the LAC curves of any two materials


$\begin{aligned} & \text { relative } \\ & \text { error }\end{aligned}=100\left|\frac{\text { Basis }- \text { Theory }}{\text { Theory }}\right|$

## Accuracy of Common 2-Element Bases v.s. Theoretical Values



## Principal Component Analysis (PCA) Basis

Dual-Energy PCA Basis


2-element PCA basis derived from LAC
values of $Z=6,7,8,13$

Triple-Energy PCA Basis


3-element PCA basis derived from LAC values of $Z=6,7,8,13,26,36$

## Accuracy of 3-Element Bases for Elements 5-36

- Dual-energy: PCA basis does not outperform conventional bases
- Triple-energy: PCA basis outperforms conventional bases


Compton, photoelectric, iron basis set


3-element PCA basis derived from LAC values of $Z=6,7,8,13,26,36$

## LAC basis plots




LAC basis here is a transformation of Compton/ photoelectric basis

$$
\begin{gathered}
\mu(E, x) \approx b_{1}(E) \mu_{1}(x)+b_{2}(E) \mu_{2}(x) \\
\mu(54, x) \approx 0 \mu_{1}(x)+1 \mu_{2}(x) \\
\mu(95, x) \approx 1 \mu_{1}(x)+0 \mu_{2}(x)
\end{gathered}
$$

## Conventional Single Energy CT

Bremsstrahlung x-ray sources provide a polychromatic spectra

Lower energies are preferentially absorbed


Polychromatic x-ray CT is a non-linear inverse problem

Conventional methods solve a linearized approximation

Leads to cupping, streaking, loss of quantitative accuracy


## Multi-Energy CT

Multi-energy CT entails measuring attenuation of two or more x-ray spectra by an object

This information may be used to remove beam hardening artifacts, improve quantitative accuracy, and enable more x-ray features to be extracted

Plot of Two X-ray Spectra


Single Energy
Reconstruction


Dual Energy Reconstruction


## Collecting Multi-Spectral CT Data

1. Scan same object twice with different spectra
2. Sandwich detectors

3. Multi-energy spectral x -ray detector energy of each detected photon is estimated

## Polychromatic CT Model




## Three or More Energies

- Triple energy decomposition results are on par
- with adaptive basis results from previous slide
- Decomposition requires solving a non-convex cost function, so converges to local minimum with a poor initialization
- Results with high Z materials (higher than 54) are variable because low energy bins have no signal
- If using a multi-spectral detector, one may be able to use adaptive binning to ensure all channels have sufficient statistics


## Approximate Polychromatic CT Model

Approximate LAC by a separable basis expansion

$$
\mu(E, x) \approx \sum_{i} b_{i}(E) c_{i}(x)
$$

Attenuation data can then be approximated by

$$
\begin{aligned}
g & \approx-\log \left(\int s(E) e^{-\sum_{i} b_{i}(E) \int_{L} c_{i}(x) d x} d E\right) \\
& =-\log \left(\int s(E) e^{-\sum_{i} b_{i}(E) a_{i}(L)} d E\right)
\end{aligned}
$$

## Multi-energy Decomposition

Suppose one has measured a collection of data, $\left(g_{1}, \ldots, g_{N}\right)$, measured with spectra $\left(\mathrm{s}_{1}, \ldots, \mathrm{~s}_{\mathrm{N}}\right)$

Then one may decompose the data into components $\left(a_{1}, \ldots, a_{N}\right)$ By minimizing the following cost function

$$
C\left(\left[\begin{array}{c}
a_{1} \\
\vdots \\
a_{N}
\end{array}\right]\right):=\frac{1}{2} \sum_{j=1}^{N}\left[g_{j}+\log \left(\int s_{j}(E) e^{-\sum_{i} b_{i}(E) a_{i}(L)} d E\right)\right]^{2}
$$

Decomposition: $\left(g_{1}, \ldots, g_{N}\right) \square\left(a_{1}, \ldots, a_{N}\right)$

## Transformation to LAC basis

Consider the Compton/ photoelectric basis $\left(b_{c}, b_{p}\right)$ and corresponding coefficients ( $\mathrm{c}_{\mathrm{c}}, \mathrm{c}_{\mathrm{p}}$ ) and energy pair $\left(\mathrm{E}_{1}, \mathrm{E}_{2}\right)$. Then

$$
\mu(E, x) \approx b_{c}(E) c_{c}(x)+b_{p}(E) c_{p}(x)
$$

We define

$$
\begin{aligned}
& \mu_{1}:=b_{c}\left(E_{1}\right) c_{c}+b_{p}\left(E_{1}\right) c_{p} \\
& \mu_{2}:=b_{c}\left(E_{2}\right) c_{c}+b_{p}\left(E_{2}\right) c_{p}
\end{aligned} \longrightarrow\left[\begin{array}{l}
\mu_{1} \\
\mu_{2}
\end{array}\right]=\left[\begin{array}{ll}
b_{c}\left(E_{1}\right) & b_{p}\left(E_{1}\right) \\
b_{c}\left(E_{2}\right) & b_{p}\left(E_{2}\right)
\end{array}\right]\left[\begin{array}{l}
c_{c} \\
c_{p}
\end{array}\right]
$$

and thus

$$
\begin{gathered}
{\left[\begin{array}{l}
c_{c} \\
c_{p}
\end{array}\right]=\left[\begin{array}{ll}
b_{c}\left(E_{1}\right) & b_{p}\left(E_{1}\right) \\
b_{c}\left(E_{2}\right) & b_{p}\left(E_{2}\right)
\end{array}\right]^{-1}\left[\begin{array}{l}
\mu_{1} \\
\mu_{2}
\end{array}\right]} \\
b_{c}(E) c_{c}(x)+b_{p}(E) c_{p}(x)=\left[b_{c}(E) b_{p}(E)\right]\left[\begin{array}{ll}
b_{c}\left(E_{1}\right) & b_{p}\left(E_{1}\right) \\
b_{c}\left(E_{2}\right) & b_{p}\left(E_{2}\right)
\end{array}\right]^{-1}\left[\begin{array}{l}
\mu_{1} \\
\mu_{2}
\end{array}\right] \\
=: b_{1}(E) \mu_{1}+b_{2}(E) \mu_{2}
\end{gathered}
$$

## LAC Basis is Just a Linear Transformation

LAC basis is simply a linear transformation of a given basis, e.g., Compton/ photoelectric, material basis, etc

If one performs multi-energy decomposition with a given basis and then transforms to an LAC basis (with a simple matrix multiplication), one get the same result by doing the multi-energy decomposition directly with an LAC basis and vice versa

No matter what the root basis used, transformation to LAC basis will give you results in LAC units

## Transforming LAC basis coefficients to Electron Density and Effective Atomic Number

One may approximate LAC by

$$
\mu(E, x) \approx \mu\left(E, Z_{e}(x)\right)=\rho_{e}(x) \sigma_{e}\left(E, Z_{e}(x)\right)
$$

where non-integer $Z_{e}$ defined by linear interpolation

If one has dual energy data one can estimate electron density and effective atomic number (effective-Z) by solving

$$
\begin{aligned}
& \mu_{1}=\rho_{e} \sigma_{e}\left(E_{1}, Z_{e}\right) \\
& \mu_{2}=\rho_{e} \sigma_{e}\left(E_{2}, Z_{e}\right)
\end{aligned}
$$

## Example of a Main Title and Content [Calibri Font]

 Use 24-pt "Regular" (no bold) subtitles to provide additional detail- Laboratory budgets over the last 15 years
- How does this affect my program?
- What are the relative values of our investments? (Discussion)
- Three critical issues to be decided:
- Size of effort
- Organization and R2A2
- Funding mechanisms
- Wrap-up

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