

Procedural Protocol for Dual Energy Computed Tomography (DECT) Inversion

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Abstract : The dual energy computed tomography (DECT) aims at noting the HU(V) values for the sample at two different voltages $V=V_1, V_2$ and thus obtain the electron densities (ρ_e) and effective atomic number (Z_{eff}) of the substance. In the present paper, we aim to obtain a numerical algorithm by which (ρ_e, Z_{eff}) can be obtained from the HU(100) and HU(140) data, where $V=100, 140$ kVp. The idea is to use this inversion method to characterize and distinguish between the lipid and fibrous coronary artery plaques. With the idea to develop the inversion algorithm for low Z_{eff} materials, as is the case with non calcified coronary artery plaque, we prepare aqueous samples whose calculated values of (ρ_e, Z_{eff}) lie in the range ($2.65 \times 10^{23} \leq \rho_e \leq 3.64 \times 10^{23}$ per cc) and ($6.80 \leq Z_{eff} \leq 8.90$). We fill the phantom with these known samples and experimentally determine HU(100) and HU(140) for the same pixels. Knowing that the HU(V) values are related to the attenuation coefficient of the system, we present an algorithm by which the (ρ_e, Z_{eff}) is calibrated with respect to (HU(100), HU(140)). The calibration is done with a known set of 20 samples; its accuracy is checked with a different set of 23 known samples. We find that the calibration gives the ρ_e with an accuracy of $\pm 4\%$ while Z_{eff} is found within $\pm 1\%$ of the actual value, the confidence being 95%. In this inversion method (ρ_e, Z_{eff}) of the scanned sample can be found by eliminating the effects of the CT machine and also by ensuring that the determination of the two unknowns (ρ_e, Z_{eff}) does not interfere with each other. It is found that this algorithm can be used for prediction of chemical characteristic (ρ_e, Z_{eff}) of unknown scanned materials with 95% confidence level, by inversion of the DECT data.

Keywords : chemical composition, dual-energy computed tomography, inversion algorithm

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