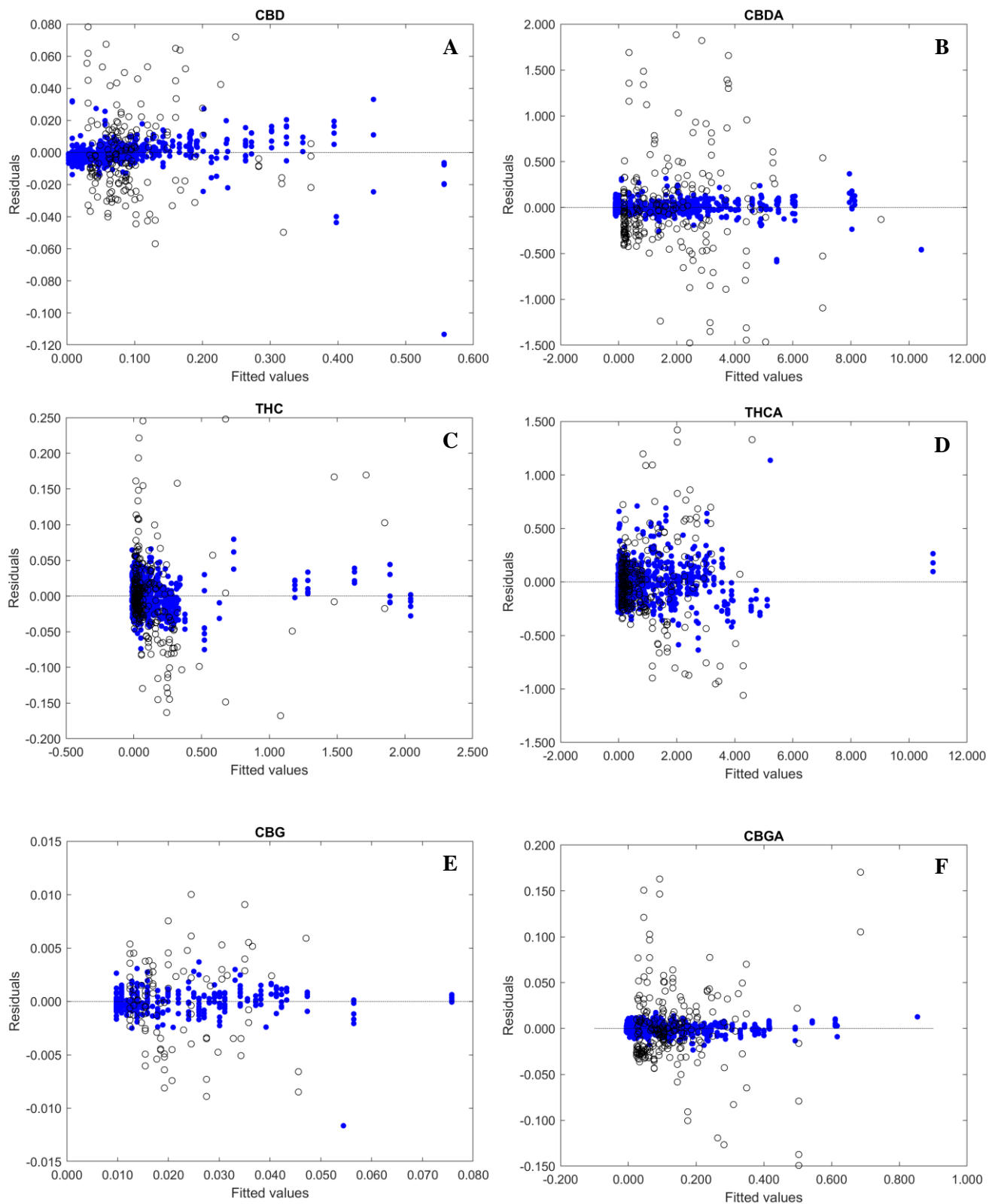


Supplementary Information Figure S1



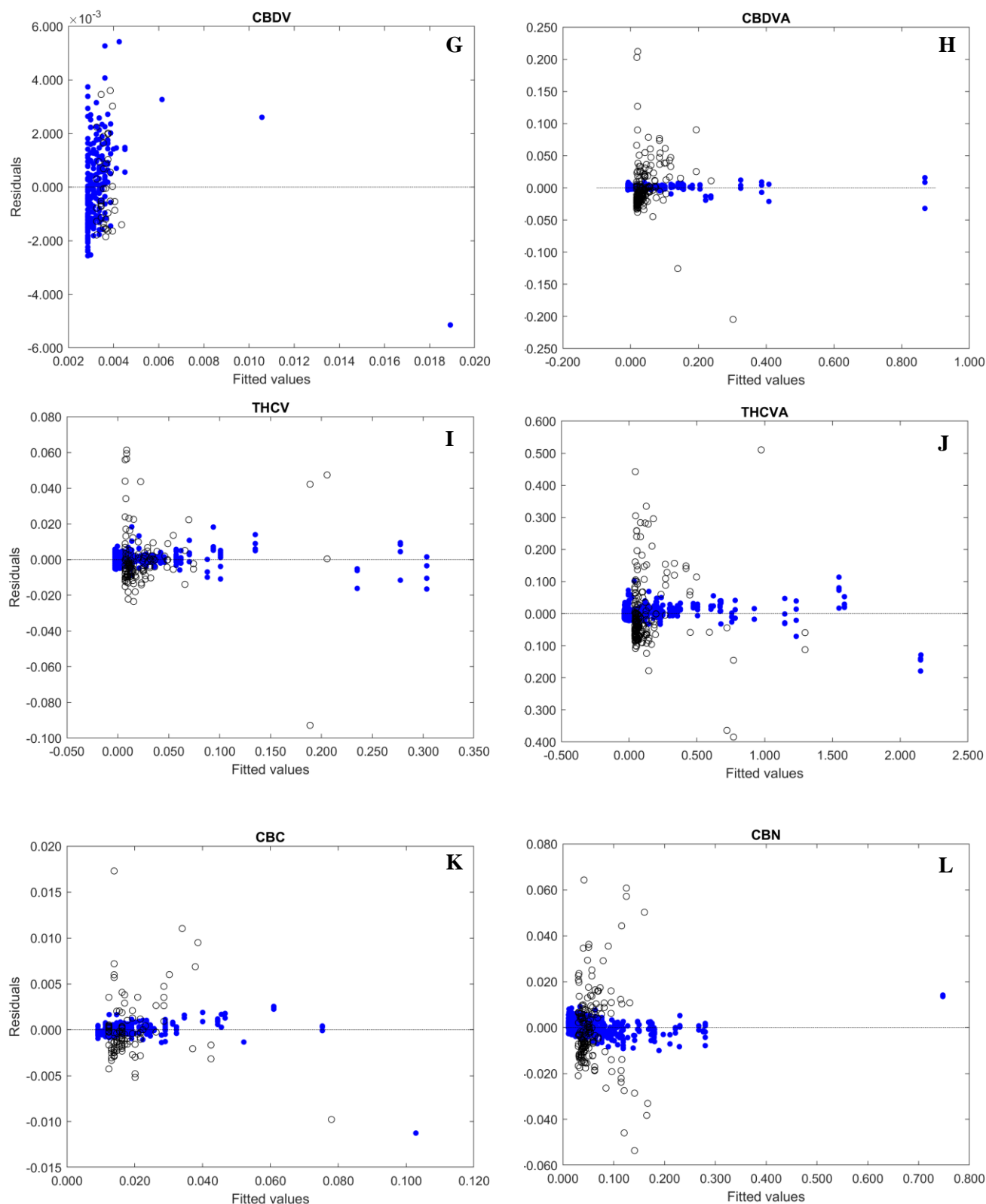


Figure S1: Residual plots for final models of all 12 target cannabinoids.

Residuals (y-axis) plotted against fitted values (x-axis) for Cross-validation (blue closed circles) and Hold-out validation (open circles) data of the 12 target Cannabinoids: A.) Cannabidiol (CBD) B.) Cannabidiolic acid (CBDA), C.) Δ^9 -tetrahydrocannabinol (Δ^9 -THC), D.) Δ^9 -tetrahydrocannabinolic acid (Δ^9 -THCA), E.) Cannabigerol (CBG), F.) Cannabigerolic acid (CBGA), G.) Cannabidivarin (CBDV), H.) Cannabidivarinic acid (CBDVA), I.) Tetrahydrocannabivarin (THCV), J.) Tetrahydrocannabivarinic acid (THCVA), K.) Cannabichromene (CBC), L.) Cannabinol (CBN).