

Webinar on

# PK/PD Studies In Drug Discovery And Development

#### **Areas Covered**

Absorption, Distribution

Metabolism, Excretion

Drug action, Dose-response relationship

Selectivity, Therapeutic window

Biomarkers, Variation in response

Hysteresis, Case study



This webinar illustrate how the model-based drug development and, in particular, PK/PD studies, can facilitate knowledge management and decision making to streamline drug discovery and development and to reduce the attrition rate.

#### **PRESENTED BY:**

Dr. Stefano Persiani is currently Director of Translational Sciences and Pharmacokinetics at Rottapharm Biotech, Italy. He graduated in Pharmacy at the University of Milan, Italy and completed a Post-Doctoral fellowship in the Department of Pathology of the University of Pittsburgh School of Medicine, Pittsburgh, Pennsylvania, USA, and later as a Research Associate in the Department of Pharmaceutics of the University of Southern California, School of Pharmacy in Los Angeles, California, USA



**On-Demand Webinar** 

**Duration: 60 Minutes** 

Price: \$200

### **Webinar Description**

The webinar reviews the general concepts and basic elements of pharmacokinetics, pharmacodynamics, and their correlation. The use of PK/PD for dose selection during drug development will be described also with the presentation of a case study. The use of PK/PD will be described for the evaluation of both drug efficacy and safety. This will illustrate how the model-based drug development and, in particular, PK/PD studies, can facilitate knowledge management and decision making to streamline drug discovery and development and to reduce the attrition rate.



#### **Who Should Attend?**

**Project Managers** 

Pre-clinical and Clinical Pharmacologists

Regulatory Affairs

Clinical Research Associates

Drug Discovery Scientists



## Why Should Attend?

PK/PD studies construct, validate, and utilize disease models, drug exposure-response models, and pharmacometrics models to facilitate drug development. PK/PD studies offer great support to learn and confirm the key characteristics of new molecular entities in a quantitative manner. This provides evidence for optimizing drug development plans and enabling critical decision-making. Drug development without PK/PD modeling is considered at a higher risk of failure.





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