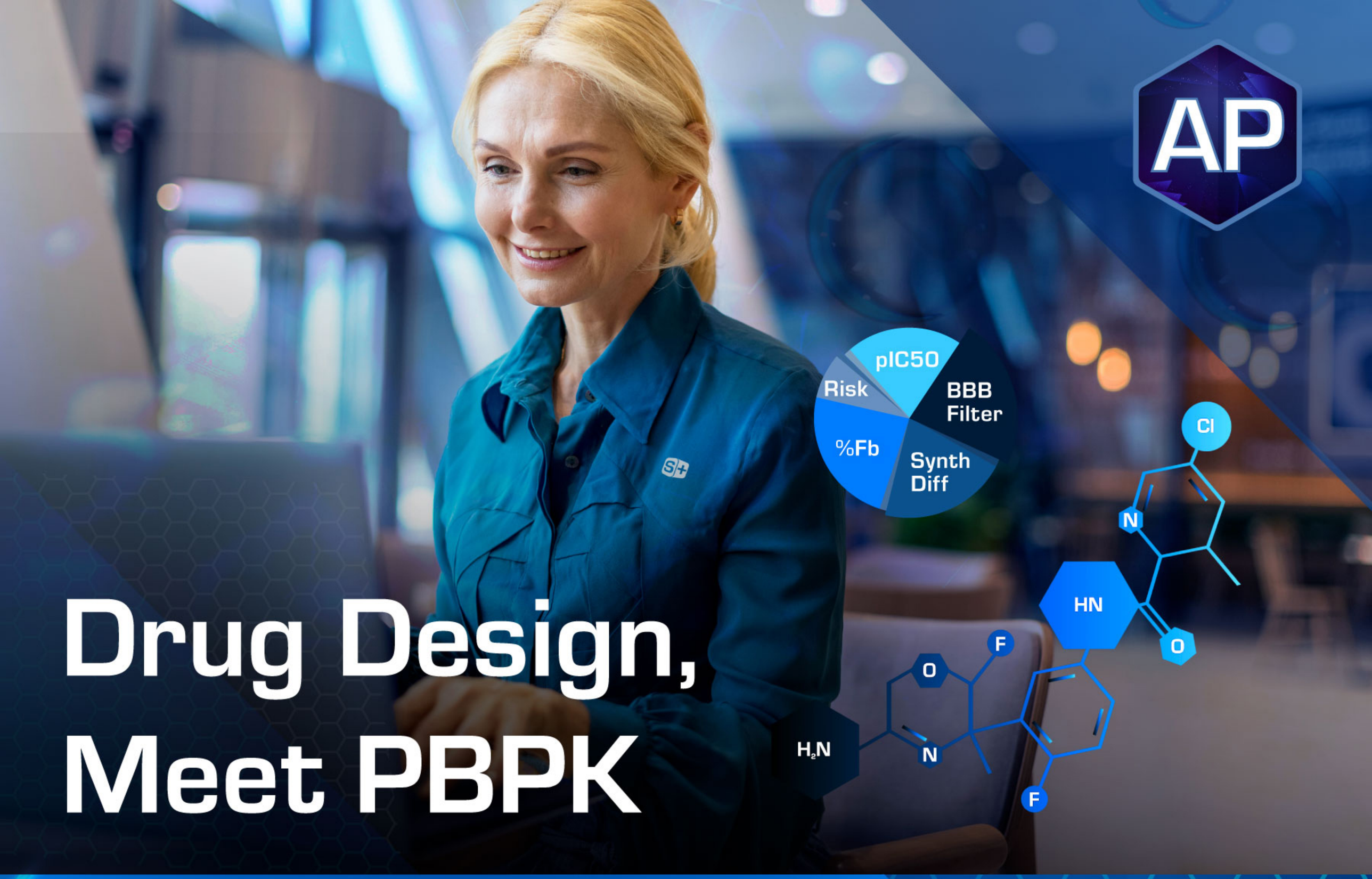

Drug Design, Meet PBPK

Leverage artificial intelligence and machine learning to automate the drug design process with industry-best ADMET and HT-PBPK predictions.

With our AI-driven Drug Design (AIDD) module, your team will be able to...

- Generate & evaluate up to 10 million molecules overnight
- Speed up hit discovery, hit-to-lead and lead optimization
- Use 3D ligand shape matching for more precise non-intuitive scaffold hopping
- Include properties from external programs like docking scores or synthetic feasibility
- Leverage Simulations Plus' 25+ years of experience in providing leading modeling & simulation software

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