

Cursos Computação UFN

Laboratório de Práticas UFN

Alexandre Zamberlan

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Computação Científica

Simulação

**MASPN** 

Sistemas Nanoparticulados Coloidais

Doutorado

Programa Nanociências 2014 a 2018

Profa. Solange Fagan

UFN

Prof. Rafael Bordini PUCRS

Ambiente de simulação integrado

Sistemas Multiagentes





Biomedicina \*

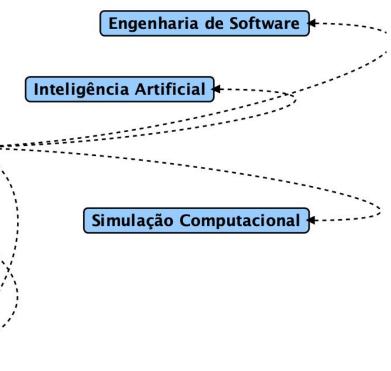
Farmácia

Física

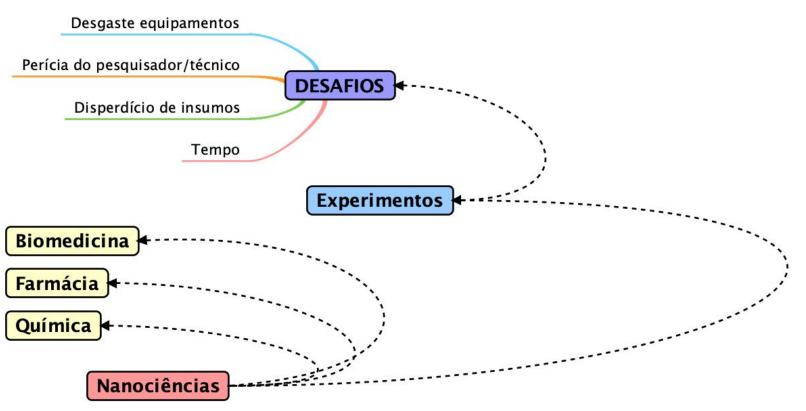
Química

Computação

Nanociências











Desgaste equipamentos

Disperdício de insumos

Tempo

Nanociências

Perícia do pesquisador/técnico

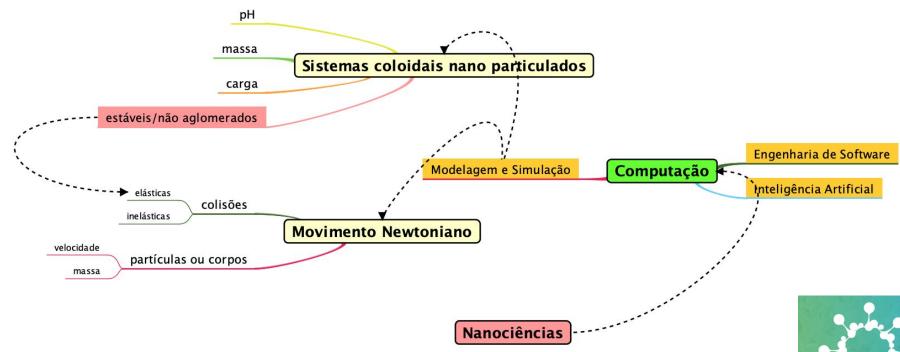
MEIO Modelagem Computação + Simulação DESAFIOS



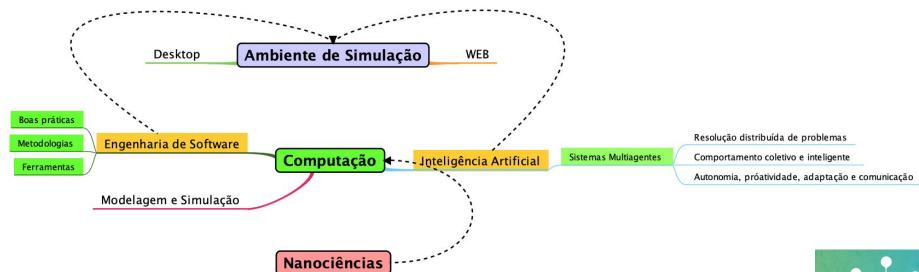
Engenharia de Software

Inteligência Artificial

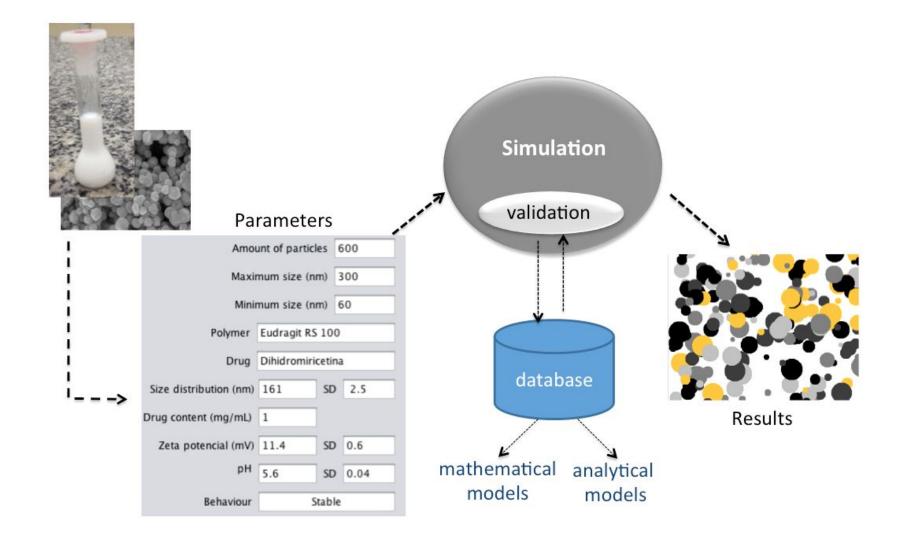












## **Y** Experiment

### H Create experiment

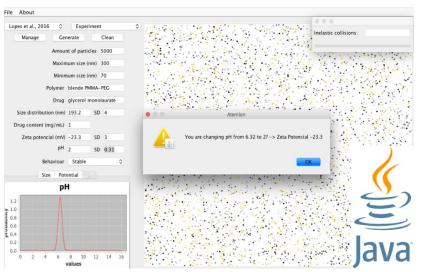
Researcher	Polymer	Drug	Drug Content (mg/mL)	Size Distribution (nm)	Size Distribution SD (%)	Zeta Potencial (mV)	Zeta Potential SD (%)	рН	pH SD (%)	Result	Edit	Delete
anajulia@unifra.br	EUDRAGIT RS 100	DMY	1.0	145.0	0.5	12.0	3.0	3.8	0.5	STABLE	Œ	•
alexz@unifra.br	EUDRAGIT RS 100	DMY	1.3	123.8	0.5	-30.1	3.0	None	None	STABLE	ß	曲
solange@unifra.br	EUDRAGIT RS 100	DMY	1.0	111.98	2.0	-12.0	2.0	6.0	1.0	NOT STABLE	Œ	
laporta@unifra.br	PCL		11111.0	112.0	1.0	<b>2</b> p	ytho	ı" ∩™	1.0	STABLE	ß	<b>*</b>
							)				0	





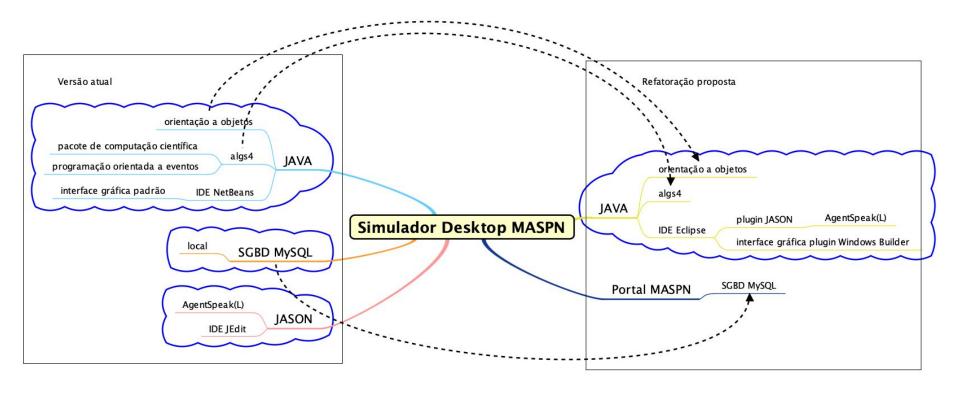












# **Y** Experiment

H Create experiment

Researcher	Polymer	Drug	Drug Content (mg/mL)	Size Distribution (nm)	Zeta Potencial (mV)	рН	D.O.I	Result	Edit	Delete
Alexandre Zamberlan	ERS100	CLOTRIMAZOL	3.0	173.0	14.4	5.6	SANTOS et al., 2014)	STABLE	Œ	
Alexandre Zamberlan	ERS100	CLOTRIMAZOL	1.0	169.0	14.5	5.7	SANTOS et al., 2014	STABLE	ß	â
Alexandre Zamberlan	ERS100	DIHIDROMIRICETINA	1.0	161.0	11.4	5.6	DALCIN et al., 2017	STABLE	(G.)	
Alexandre Zamberlan	ERS100	DIHIDROMIRICETINA	5.0	123.0	13.4	3.8	DALCIN et al., 2017	NOT STABLE	©	
Alexandre Zamberlan	ERS100	DIHIDROMIRICETINA	2.0	151.0	12.7	4.2	DALCIN et al., 2017	STABLE	©	â
Alexandre Zamberlan	PCL	NISINA	1.0	234.0	-6.62	5.3	ABREU et al., 2016	STABLE	Œ	â





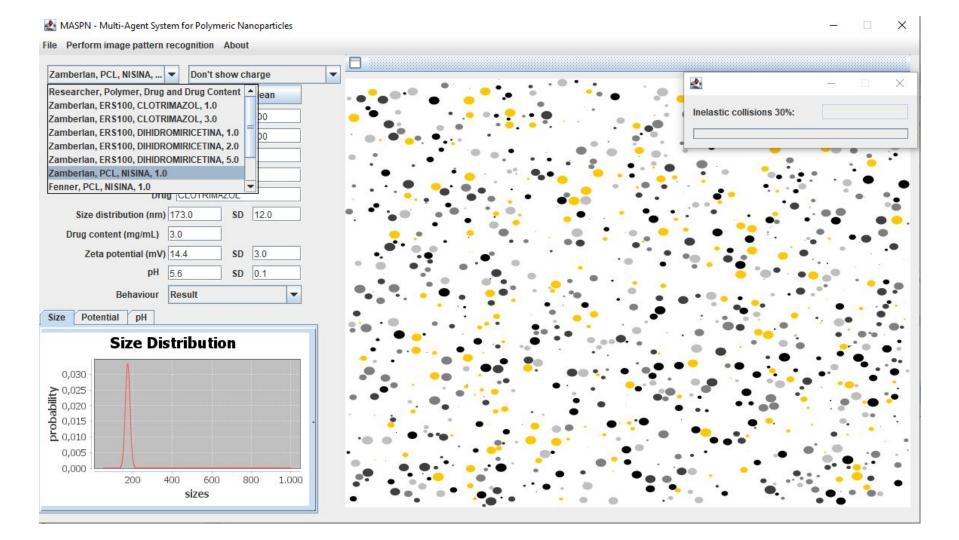


H Create polymer

Initiale

Description	Initials	Description	Initials	
Anfotericina B	ANFOTERICINA B	Eudragit RL100	ERL100	
Clotrimazol	CLOTRIMAZOL	Eudragit RS100	ERS100	
Curcumina	CURCUMINA	Polycaprolactone	PCL	
Dihidromiricetina	DIHIDROMIRICETINA	Polylactic acid	PLA	
Nisina	NISINA	PLA + Pluronic F68	PLA + PLURONIC F68	
Pravastiatina	PRAVASTIATINA	PLA + Solutol HS15	PLA + SOLUTOL HS15	
Sparfloxacino	SPARFLOXACINO	Poly(lactic-co-glycolic acid)	PLGA	
Tretinoína	TRETINOÍNA	Quitosana	QUITOSANA	

Description







# Portal of experiments on polymer nanoparticles for simulation in MASPN

The MASPN simulation tool has database access with experiments published in scientific papers with production data and characterization of polymeric nanoparticles (PNPs). Thus, a collaborative portal among institutions conducting research on PNPs assists in the discovery of knowledge of best practices in nanoparticle production.

### Some published results:

- · A simulation environment for polymeric nanoparticles based on multi-agent systems
- Multi-Agent Systems, Simulation and Nanotechnology
- Sistema Multiagente para avaliação do efeito de aglomeração em partículas poliméricas

This project is associeted to the Laboratory of Practices of the undergraduate courses in Computer Science and Information Systems and the LASIMON at Universidade Franciscana.



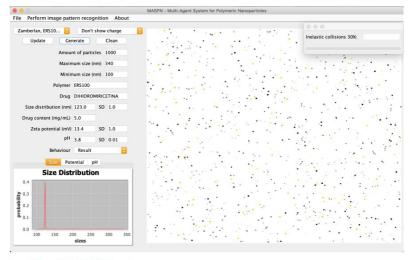






# Download

You can download MASPN tool as compressed file.



### ATTENTION! There are details for running the MASPN tool.

### Windows Users:

- Java JDK installed
- Double click the maspnV2.jar icon

### Linux and Mac OS Users:

- Java JDK installed
- Run in shell the command "java -jar maspnV2.jar"



# About

The MASPN simulation tool has database access with experiments published in scientific papers with production data and characterization of polymeric nanoparticles (PNPs). Thus, a collaborative portal among institutions conducting research on PNPs assists in the discovery of knowledge of best practices in nanoparticle production.

### Collaborators:

#### UFN

- Alexandre Zamberlan Computer science researcher
- o Ana Júlia Dalcin Pharmacy researcher
- o Ederson Guterres Costa Information System undergraduate student
- o Guilherme Chagas Kurtz Computer science researcher
- o Gustavo Pereira Computer Science undergraduate student
- o Lúcio Colusso Barnewitz Computer Science undergraduate student
- o Pierre Carvalho Fenner Computer Science undergraduate student
- Solange Fagan Physics researcher
- o Sylvio Vieira Computer science researcher
- o Tomás Gomes Loureiro Computer Science undergraduate student

#### PUCRS

o Rafael Heitor Bordini - Computer science researcher

### FEEVALE

- Fernando Dal Pont Morisso Chemistry researcher
- o Kevin Brandon Villa Chemical Engineering undergraduate student

#### • ER Clinic

Robertson Ebling dos Santos - Information Technology professional









### REPÚBLICA FEDERATIVA DO BRASIL

MINISTÉRIO DA ECONOMIA

### INSTITUTO NACIONAL DA PROPRIEDADE INDUSTRIAL

DIRETORIA DE PATENTES, PROGRAMAS DE COMPUTADOR E TOPOGRAFIAS DE CIRCUITOS INTEGRADOS

# Certificado de Registro de Programa de Computador

Processo No: BR512021001965-2

O Instituto Nacional da Propriedade Industrial expede o presente certificado de registro de programa de computador, válido por 50 anos a partir de 1° de janeiro subsequente à data de 01/07/2021, em conformidade com o \$2°, art. 2° da Lei 9.609, de 19 de Fevereiro de 1998.

Título: MASPN - Multi-agent System for Polymeric Nanoparticle

Data de publicação: 01/07/2021

Data de criação: 04/04/2020

Titular(es): SOCIEDADE CARITATIVA E LITERÁRIA SÃO FRANCISCO DE ASSIS ZONA NORTE

Autor(es): ALEXANDRE DE OLIVEIRA ZAMBERLAN; ROBERTSON EBLING DOS SANTOS; SOLANGE BINOTTO FAGAN; TOMÁS LOUREIRO GOMES; GUILHERME CHAGAS KURTZ; PIERRE CARVALHO FENNER; SYLVIO ANDRÉ GARCIA VIEIRA; ANA JÚLIA FIGUEIRÓ DALCIN; ÉDERSON GUTERRES COSTA; GUSTAVO GARCIA PEREIRA; LUCIO COLUSSO BARNEWITZ; RAFAEL HEITOR BORDINI



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**MASPN** 

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Validação do portal: fármacos e polímeros

Mineração de Dados: descoberta de padrões

Computação Paralela: GPU

**SMA Reativos versus Cognitivos** 

Validação do simulador com novos fármacos e polímeros

Adequações do ambiente ao grupo de pesquisa

