Resultados de prueba de plagio: org-109.txt

Plagio detectado: 90.32%

Texto original: of research for pharmaceutical companies and chemical scientists.
Texto plagiado: of research for pharmaceutical companies and chemical scientists.
Texto original: rug design
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Texto plagiado: rug design
Texto original: Drug designing and development is an important area of research for
pharmaceutical companies and chemical scientists.
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Texto original: low efficacy, off-target delivery, time consumption, and high cost

Texto plagiado: low efficacy, off-target delivery, time consumption, and high cost

Texto original: However, low efficacy, off-target delivery, time consumption, and high cost impose a hurdle and challenges that impact drug design and discovery.

Texto plagiado: However, low efficacy, off-target delivery, time consumption, and high cost impose a hurdle and challenges that impact drug design and discovery.

Texto original: drug design and discovery

Texto plagiado: drug design and discovery

Texto original: data from genomics, proteomics, microarray data, and clinical trials

Texto plagiado: data from genomics, proteomics, microarray data, and clinical trials

Texto original: drug discovery

Texto plagiado: drug discovery

Texto original: Further, complex and big data from genomics, proteomics, microarray data, and clinical trials also impose an obstacle in the drug discovery pipeline.

Texto plagiado: Further, complex and big data from genomics, proteomics, microarray data,

and clinical trials also impose an obstacle in the drug discovery pipeline.
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Texto original: Artificial intelligence and machine learning technology play a crucial role in drug discovery and development.

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Texto original: peptide synthesis, structure-based virtual screening, ligand-based virtual screening, toxicity prediction, drug monitoring and release, pharmacophore modeling, quantitative structure–activity relationship, drug repositioning, polypharmacology, and physiochemical activity.

Texto plagiado: peptide synthesis, structure-based virtual screening, ligand-based virtual screening, toxicity prediction, drug monitoring and release, pharmacophore modeling, quantitative structure–activity relationship, drug repositioning, polypharmacology, and

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Texto original: Machine learning and deep learning algorithms have been implemented in several drug discovery processes such as peptide synthesis, structure-based virtual screening, ligand-based virtual screening, toxicity prediction, drug monitoring and release, pharmacophore modeling, quantitative structure–activity relationship, drug repositioning,

polypharmacology, and physiochemical activity.

Texto plagiado: Machine learning and deep learning algorithms have been implemented in several drug discovery processes such as peptide synthesis, structure-based virtual screening, ligand-based virtual screening, toxicity prediction, drug monitoring and release, pharmacophore modeling, quantitative structure–activity relationship, drug repositioning, polypharmacology, and physiochemical activity.

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Texto original: artificial intelligence and deep learning

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Texto original: , novel data mining, curation, and management techniques

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