

1. Physicochemical Property

Property	Value	Comment
Molecular Weight	346.21	Contain hydrogen atoms. Optimal:100~600
Volume	370.718	Van der Waals volume
Density	0.934	Density = MW / Volume
nHA	4	Number of hydrogen bond acceptors. Optimal:0~12
nHD	2	Number of hydrogen bond donors. Optimal:0~7
nRot	3	Number of rotatable bonds. Optimal:0~11
nRing	3	Number of rings. Optimal:0~6
MaxRing	14	Number of atoms in the biggest ring. Optimal:0~18
nHet	4	Number of heteroatoms. Optimal:1~15
fChar	0	Formal charge. Optimal:-4 ~4
nRig	17	Number of rigid bonds. Optimal:0~30
Flexibility	0.176	Flexibility = nRot /nRig
Stereo Centers	2	Optimal: ≤ 2
TPSA	66.76	Topological Polar Surface Area. Optimal:0~140
logS	-4.35	Log of the aqueous solubility. Optimal: -4~0.5 log mol/L
logP	4.544	Log of the octanol/water partition coefficient. Optimal: 0~3
logD	3.778	logP at physiological pH 7.4. Optimal: 1~3

2. Medicinal Chemistry

Property	Value	Decision	Comment
QED	0.615	•	■ A measure of drug-likeness based on the concept of desirability; ■ Attractive: > 0.67; unattractive: 0.49~0.67; too complex: < 0.34
SAscore	3.779	•	 ■ Synthetic accessibility score is designed to estimate ease of synthesis of drug-like molecules. ■ SAscore ≥ 6, difficult to synthesize; SAscore <6, easy to synthesize
Fsp3	0.667	•	 ■ The number of sp3 hybridized carbons / total carbon count, correlating with melting point and solubility. ■ Fsp³ ≥0.42 is considered a suitable value.
MCE-18	79.514	•	■ MCE-18 stands for medicinal chemistry evolution.■ MCE-18≥45 is considered a suitable value.

NPscore	2.129	-	■ Natural product-likeness score. ■ This score is typically in the range from –5 to 5. The higher the score is, the higher the probability is that the molecule is a NP.
Lipinski Rule	Accepted	•	 ■ MW ≤ 500; logP ≤ 5; Hacc ≤ 10; Hdon ≤ 5 ■ If two properties are out of range, a poor absorption or permeability is possible, one is acceptable.
Pfizer Rule	Rejected	•	logP > 3; TPSA < 75 Compounds with a high log P (>3) and low TPSA (<75) are likely to be toxic.
GSK Rule	Rejected	•	 ■ MW ≤ 400; logP ≤ 4 ■ Compounds satisfying the GSK rule may have a more favorable ADMET profile
Golden Triangle	Accepted	•	 ■ 200 ≤ MW ≤ 50; -2 ≤ logD ≤ 5 ■ Compounds satisfying the Golden Triangle rule may have a more favorable ADMET profile.
PAINS	1 alerts	-	Pan Assay Interference Compounds, frequent hitters, Alpha-screen artifacts and reactive compound.
ALARM NMR	1 alerts	-	Thiol reactive compounds.
BMS	0 alerts	-	Undesirable, reactive compounds.
Chelator Rule	1 alerts	-	Chelating compounds.

3. Absorption

Property	Value	Decision	Comment
Caco-2 Permeability	-4.735	•	Optimal: higher than -5.15 Log unit
MDCK Permeability	1.9e-05	•	 ■ low permeability: < 2 x 10⁻⁶ cm/s ■ medium permeability: 2-20 x 10⁻⁶ cm/s ■ high passive permeability: > 20 x 10⁻⁶ cm/s
Pgp-inhibitor	0.757	•	■ Category 1: Inhibitor; Category 0: Non-inhibitor; ■ The output value is the probability of being Pgp-inhibitor
Pgp-substrate	0.001	•	■ Category 1: substrate; Category 0: Non-substrate; ■ The output value is the probability of being Pgp-substrate
НІА	0.014	•	■ Human Intestinal Absorption ■ Category 1: HIA+(HIA < 30%); Category 0: HIA-(HIA < 30%); The output value is the probability of being HIA+
F _{20%}	0.989	•	■ 20% Bioavailability ■ Category 1: $F_{20\%}$ + (bioavailability < 20%); Category 0: $F_{20\%}$ - (bioavailability ≥ 20%); The output value is the probability of being $F_{20\%}$ +

F _{30%}	0.965	•	■ 30% Bioavailability ■ Category 1: $F_{30\%}$ + (bioavailability < 30%); Category 0: $F_{30\%}$ - (bioavailability ≥ 30%); The output value is the probability of being $F_{30\%}$ +
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4. Distribution

Property	Value	Decision	Comment
PPB	98.85%	•	■ Plasma Protein Binding■ Optimal: < 90%. Drugs with high protein-bound may have a low therapeutic index.
VD	2.416	•	■ Volume Distribution ■ Optimal: 0.04-20L/kg
BBB Penetration	0.487	•	■ Blood-Brain Barrier Penetration ■ Category 1: BBB+; Category 0: BBB-; The output value is the probability of being BBB+
Fu	1.655%	•	■ The fraction unbound in plasms■ Low: <5%; Middle: 5~20%; High: > 20%

5. Metabolism

Property	Value	Comment
CYP1A2 inhibitor	0.204	■ Category 1: Inhibitor; Category 0: Non-inhibitor;■ The output value is the probability of being inhibitor.
CYP1A2 substrate	0.925	■ Category 1: Substrate; Category 0: Non-substrate;■ The output value is the probability of being substrate.
CYP2C19 inhibitor	0.449	■ Category 1: Inhibitor; Category 0: Non-inhibitor;■ The output value is the probability of being inhibitor.
CYP2C19 substrate	0.875	■ Category 1: Substrate; Category 0: Non-substrate;■ The output value is the probability of being substrate.
CYP2C9 inhibitor	0.421	■ Category 1: Inhibitor; Category 0: Non-inhibitor; ■ The output value is the probability of being inhibitor.
CYP2C9 substrate	0.927	■ Category 1: Substrate; Category 0: Non-substrate; ■ The output value is the probability of being substrate.
CYP2D6 inhibitor	0.306	■ Category 1: Inhibitor; Category 0: Non-inhibitor; ■ The output value is the probability of being inhibitor.
CYP2D6 substrate	0.806	■ Category 1: Substrate; Category 0: Non-substrate;■ The output value is the probability of being substrate.
CYP3A4 inhibitor	0.359	■ Category 1: Inhibitor; Category 0: Non-inhibitor; ■ The output value is the probability of being inhibitor.
CYP3A4 substrate	0.434	■ Category 1: Substrate; Category 0: Non-substrate; ■ The output value is the probability of being substrate.

6. Excretion

Property	Value	Decision	Comment
CL	4.21	•	■ Clearance ■ High: >15 mL/min/kg; moderate: 5-15 mL/min/kg; low: <5 mL/min/kg
T _{1/2}	0.238	-	 ■ Category 1: long half-life; Category 0: short half-life; ■ long half-life: >3h; short half-life: <3h ■ The output value is the probability of having long half-life.

7. Toxicity

Property	Value	Decision	Comment	
hERG Blockers	0.006	•	■ Category 1: active; Category 0: inactive;■ The output value is the probability of being active.	
н-нт	0.169	•	 ■ Human Hepatotoxicity ■ Category 1: H-HT positive(+); Category 0: H-HT negative(-); ■ The output value is the probability of being toxic. 	
DILI	0.027	•	 ■ Drug Induced Liver Injury. ■ Category 1: drugs with a high risk of DILI; Category 0: drugs with no risk of DILI. The output value is the probability of being toxic. 	
AMES Toxicity	0.043	•	■ Category 1: Ames positive(+); Category 0: Ames negative(-);■ The output value is the probability of being toxic.	
Rat Oral Acute Toxicity	0.439	•	 ■ Category 0: low-toxicity; Category 1: high-toxicity; ■ The output value is the probability of being highly toxic. 	
FDAMDD	0.735	•	 ■ Maximum Recommended Daily Dose ■ Category 1: FDAMDD (+); Category 0: FDAMDD (-) ■ The output value is the probability of being positive. 	
Skin Sensiti zation	0.919	•	■ Category 1: Sensitizer; Category 0: Non-sensitizer; ■ The output value is the probability of being sensitizer.	
Carcinogen city	0.061	•	■ Category 1: carcinogens; Category 0: non-carcinogens;■ The output value is the probability of being toxic.	
Eye Corrosion	0.003	•	■ Category 1: corrosives ; Category 0: noncorrosives ■ The output value is the probability of being corrosives.	
Eye Irritation	0.834	•	■ Category 1: irritants; Category 0: nonirritants ■ The output value is the probability of being irritants.	

Respiratory Toxicity	0.724	•	■ Category 1: respiratory toxicants; Category 0: respiratory nontoxicants ■ The output value is the probability of being toxic.
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8. Environmental toxicity

Property	Value	Comment
Bioconcentration Factors	1.837	■ Bioconcentration factors are used for considering secondary poisoning potential and assessing risks to human health via the food chain. ■ The unit is -log10[(mg/L)/(1000*MW)]
IGC ₅₀	4.71	 ■ Tetrahymena pyriformis 50 percent growth inhibition concentration ■ The unit is -log10[(mg/L)/(1000*MW)]
LC ₅₀ FM	4.862	■ 96-hour fathead minnow 50 percent lethal concentration ■ The unit is -log10[(mg/L)/(1000*MW)]
LC ₅₀ DM	6.225	■ 48-hour daphnia magna 50 percent lethal concentration ■ The unit is -log10[(mg/L)/(1000*MW)]

9. Tox21 pathway

Property	Value	Decision	Comment
NR-AR	0.233	•	 ■ Androgen receptor ■ Category 1: actives; Category 0: inactives; ■ The output value is the probability of being active.
NR-AR-LBD	0.033	•	 ■ Androgen receptor ligand-binding domain ■ Category 1: actives; Category 0: inactives; ■ The output value is the probability of being active.
NR-AhR	0.224	•	 ■ Aryl hydrocarbon receptor ■ Category 1: actives; Category 0: inactives; ■ The output value is the probability of being active.
NR-Aromatase	0.836	•	■ Category 1: actives; Category 0: inactives;■ The output value is the probability of being active.
NR-ER	0.585	•	 ■ Estrogen receptor ■ Category 1: actives; Category 0: inactives; ■ The output value is the probability of being active.
NR-ER-LBD	0.731	•	 ■ Estrogen receptor ligand-binding domain ■ Category 1: actives; Category 0: inactives; ■ The output value is the probability of being active.
NR-PPAR- gamma	0.963	•	 ■ Peroxisome proliferator-activated receptor gamma ■ Category 1: actives; Category 0: inactives; ■ The output value is the probability of being active.
SR-ARE	0.8	•	 ■ Antioxidant response element ■ Category 1: actives; Category 0: inactives; ■ The output value is the probability of being active.
SR-ATAD5	0.377	•	■ ATPase family AAA domain-containing protein 5 ■ Category 1: actives; Category 0: inactives; ■ The output value is the probability of being active.

SR-HSE	0.899	•	 ■ Heat shock factor response element ■ Category 1: actives; Category 0: inactives; ■ The output value is the probability of being active.
SR-MMP	0.932	•	 ■ Mitochondrial membrane potential ■ Category 1: actives; Category 0: inactives; ■ The output value is the probability of being active.
SR-p53	0.952	•	■ Category 1: actives; Category 0: inactives; ■ The output value is the probability of being active.

10. Toxicophore Rules

Property	Value	Comment
Acute Toxicity Rule	0 alerts	■ 20 substructures■ acute toxicity during oral administration
Genotoxic Carcinogenicity Rule	0 alerts	■ 117 substructures ■ carcinogenicity or mutagenicity
NonGenotoxic Carcinogenicity Rule	0 alerts	■ 23 substructures ■ carcinogenicity through nongenotoxic mechanisms
Skin Sensitization Rule	5 alerts	■ 155 substructures ■ skin irritation
Aquatic Toxicity Rule	0 alerts	■ 99 substructures ■ toxicity to liquid(water)
NonBiodegradable Rule	1 alerts	■ 19 substructures ■ non-biodegradable
SureChEMBL Rule	1 alerts	■ 164 substructures ■ MedChem unfriendly status