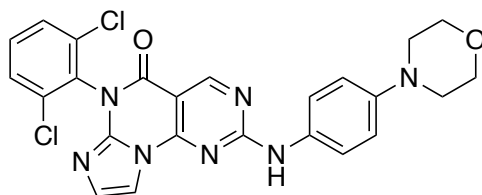


WEE2



AKI00000020a

Chemical Name: 6-(2,6-dichlorophenyl)-2-((4-morpholinophenyl)amino)imidazo[1,2-a]pyrimido[5,4-e]pyrimidin-5(6H)-one

CHEBI: 39790

Smile String:

O=C(N(C1=C(Cl)C=CC=C1Cl)C2=NC=CN23)C4=C3N=C(NC5=CC=C(N6CCOC6)C=C5)N=C4

Chemical Formula: C₂₄H₁₉Cl₂N₇O₂

Molecular Weight: 508.36

cLogP: 1.803

Source: SGC-UNC (Abbvie donation)

Reference: N/A

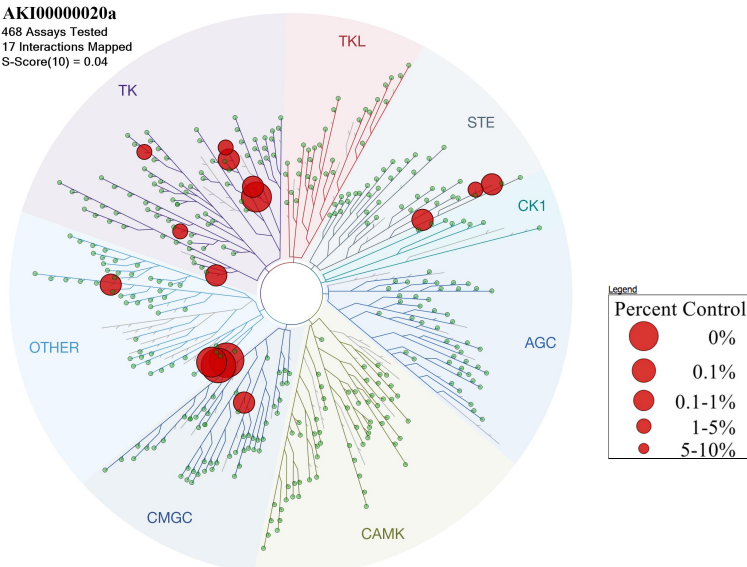
Biochemical profiling

DiscoverX (403 wild-type human kinases)

S₁₀ (1μM): 0.037 (15 kinase < 10% control)

WEE2 IC₅₀ = 10 nM

AKI00000020a
468 Assays Tested
17 Interactions Mapped
S-Score(10) = 0.04



Kinase	% Control @ 1uM
WEE1	0.1
WEE2	0.1
DDR1	0.2
EIF2AK4	0.9
DDR2	1.4
NEK2	1.6
MAP3K3	1.9
MAP3K4	2
MAPK6	2.6
SYK	3.3
JAK2	4.3
LCK	5.5
MAP3K2	6.3
JAK3	7.1
CSF1R	9.6

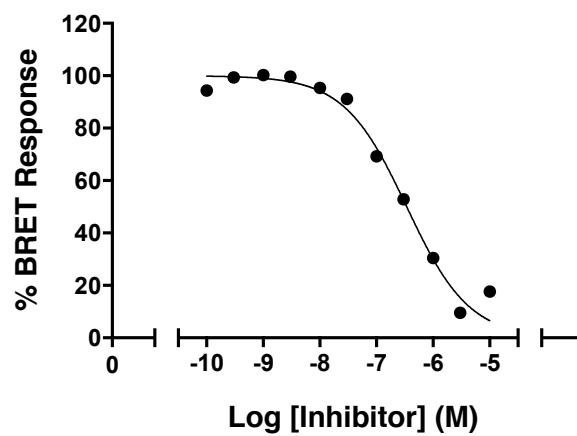
a. Treespot of DiscoverX KINOMEScan data. b. List of kinases inhibited < 10% control

Cellular target engagement in HEK293 cells

NLuc-WEE2 (N term)

WEE2 IC₅₀ = 342 nM

AKI00000020a



Cellular target engagement of AKI00000020a with WEE2