



















Structure	SMILES	BIPAD name	Name	MW	Plate description	QAS	Outlog number	Concentration	Compound description	Solvent	Plate part number
	<chem>COCC1=CC(O)=C(O)C(O)C1=C1C=CC(OC)C(O)C(O)C1O</chem>	5-(methoxymethyl)-4-[(2,3,4-trihydroxy-6-(methoxymethyl)phenyl)benzene-1,2,3-triol]	2,2',3,3',4,4'-Hexahydroxy-1,1'-biphenyl-6,6'-dimethandimethyl ether	338.3	Kinase Inhibitor Library		EI-273	10nM	PKC alpha, PKC gamma	DMSO	2832
	<chem>O=C1C2=CC=CC=C2C3=CC=CC=C3C4=CC=CC=C4C5=CC=CC=C51</chem>	14,15-diazatetracyclo[7.8.10^0.0^0]^1,1'-hexadeca-1(15),2,4,6,8,16,10,12-heptam-9-one	SP 600125	292.2	Kinase Inhibitor Library		129-58-6	EI-305	JNK	DMSO	2832
	<chem>O=C1NC2=CC=CC=C2C(=O)C1=CN3C=CC=CC3=O</chem>	2-(2-oxo-2,3-dihydro-1H-indol-3-ylidene)-2,3-dihydro-1H-indol-3-one	Indirubin	262.3	Kinase Inhibitor Library		479-41-4	CO-206	GSK-3beta, CDK5	DMSO	2832
	<chem>O=C1C=CC(=O)C=C1C(=O)C2=CC=CC=C2C3=CC=CC=C3</chem>	3-[(2E)-3-(hydroxymino)-2,3-dihydro-1H-indol-2-ylidene]-2,3-dihydro-1H-indol-2-one	Indirubin-3'-monooxime	277.3	Kinase Inhibitor Library		100807-49-8	CO-207	GSK-3beta	DMSO	2832
	<chem>ClC1=CC=CC=C1C(=O)N[C@@H](C)C1=CC=CC=C1</chem>	4-[(1S)-1-aminoethyl]-N-(pyridin-4-yl)cyclohexane-1-carboxamide dihydrochloride	Y-27632 (HCl)	320.3	Kinase Inhibitor Library		148988-50-7	EI-299	ROCK	DMSO	2832
	<chem>BrC1=CC2=CC=CC=C2C3=CC=CC=C3C4=CC=CC=C4C5=CC=CC=C51</chem>	14-bromo-8,18-diazatetracyclo[9.7.0.0^0.0^0]^1,1'-octadeca-1(11),2(7),3,5,12(17),13,15-heptam-9-one	Kempulone	327.2	Kinase Inhibitor Library		142373-20-9	EI-310	GSK-3beta	DMSO	2832
	<chem>O=C1C=CC(=O)C=C1C</chem>	1R,6S)-3-hydroxy-4-methyl-7-oxabicyclo[4.1.0]hept-3-ene-2,5-dione	Tartric acid	154.1	Kinase Inhibitor Library		121-40-4	EI-328	BTK	DMSO	2832
	<chem>OC1=CC=CC=C1C(=O)N[C@@H](C)C1=CC=CC=C1</chem>	2R,3R,4S,5R)-2-(5-amino-7-methyl-2,6,7,8,11-pentazatricyclo[3.3.1.0^1.0^1]^3dodeca-1(12),3,5,8,10-pentam-2-yl)-5-hydroxymethylisoleane-3,4-diol	Tricirbine	320.3	Kinase Inhibitor Library		35943-35-2	EI-332	Akt signaling pathway	DMSO	2832
	<chem>OC(=O)C1=CC=CC=C1C2=CC=CC=C2C3=CC=CC=C3</chem>	2,3-diphenylisoindolin-9-carboxylic acid	BML-257	326.3	Kinase Inhibitor Library		32387-96-5	EI-338	Akt	DMSO	2832
	<chem>NC(=O)C1=CC=CC=C1C2=CC=CC=C2</chem>	3-amino-5-(thiophen-3-yl)thiophene-2-carboxamide	SO-514	224.3	Kinase Inhibitor Library		354812-17-2	EI-343	IKK2	DMSO	2832
	<chem>CC(=O)N(C)C(=O)C1=CC=CC=C1</chem>	2-phenyl-N-[5-(propan-2-yl)-1,3-thiazol-2-yl]acetamide	BML-259	260.4	Kinase Inhibitor Library		207654-00-2	EI-344	Cdk5/p25	DMSO	2832
	<chem>O=C1C=CC(=O)C=C1C(=O)C2=CC=CC=C2</chem>	5,7-dihydroxy-2-(4-hydroxyphenyl)-6H-chromen-4-one	Ajigenin	270.2	Kinase Inhibitor Library		620-36-5	EI-345	CK-1B	DMSO	2832
	<chem>CCOC1=CC=CC=C1C2=CC=CC=C2</chem>	N-(3-ethoxyphenyl)-6,7-dimethylquinazolin-4-amine	BML-285	305.3	Kinase Inhibitor Library		EI-346	10nM	EGFRK	DMSO	2832
	<chem>OC1=CC=CC=C1C(=O)N[C@@H](C)C1=CC=CC=C1</chem>	1R,8S,12S,15R,16R,18R,19R,21R,23S,24E,26E,28E,30S,32S,35R)-1,8-dihydroxy-12-[[2R)-1-[(1S,3R,4R)-4-hydroxy-3-methoxycyclohexyl]propan-2-yl]-18,30-dimethoxy-15,17,21,23,25,29,35-hexamethyl-11,38-dioxo-4-azatricyclo[20.3.1.0^1.0^1]^1,1'-hexatriaconta-16,24,26,28-tetraene-2,3,10,14,20-pentone	Rapamycin	914.2	Kinase Inhibitor Library		53123-88-9	A-275	mTOR	DMSO	2832