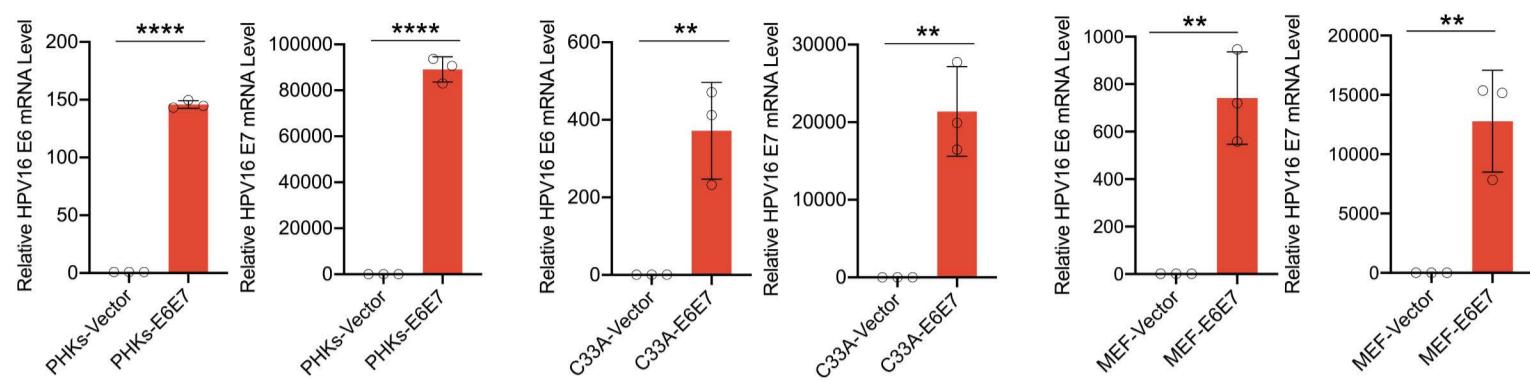
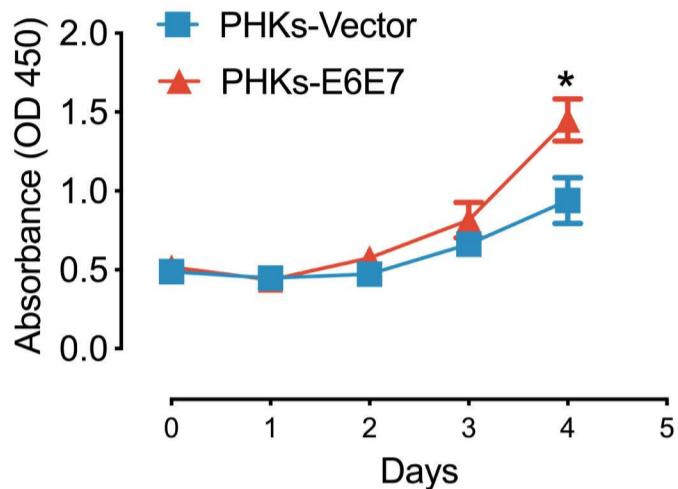


Supplementary Figure 1

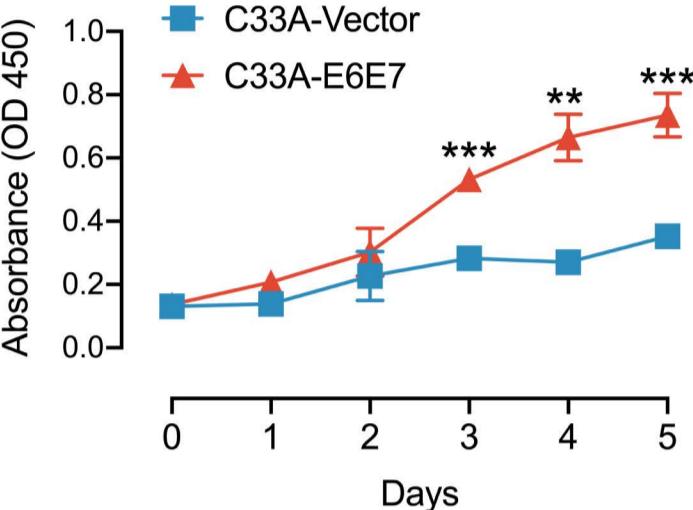
S1A



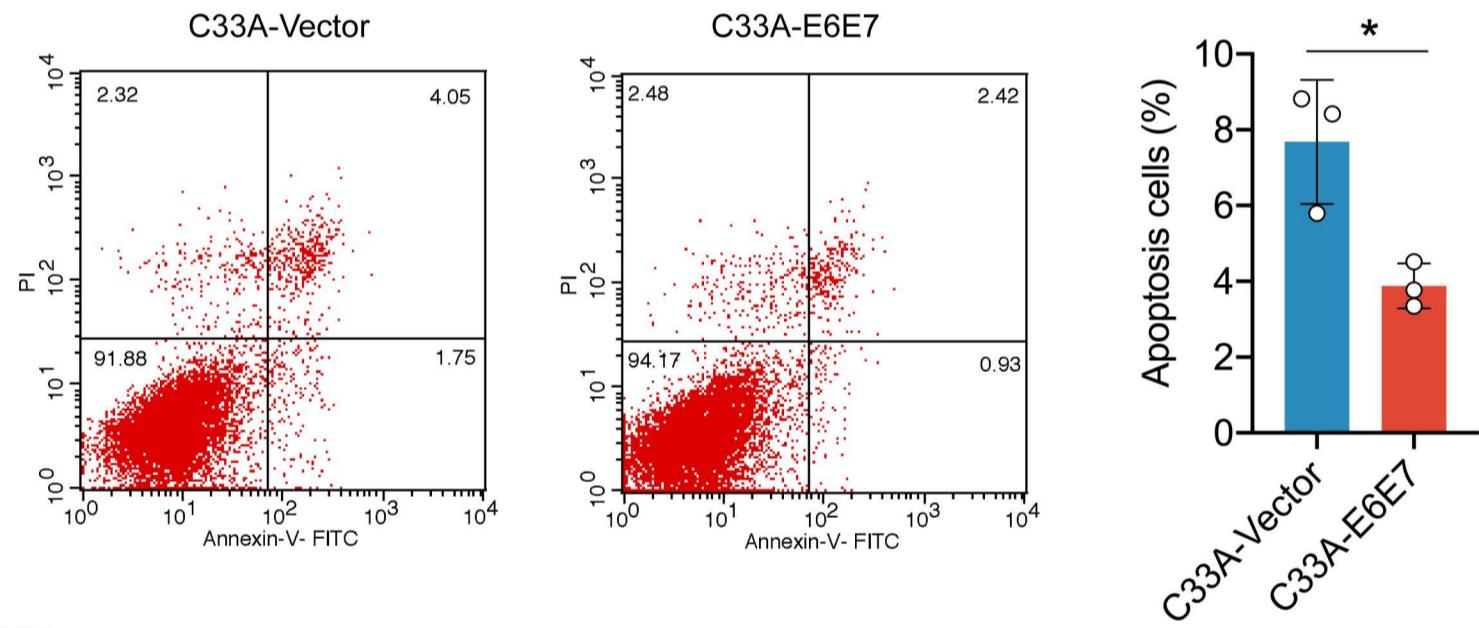
S1B



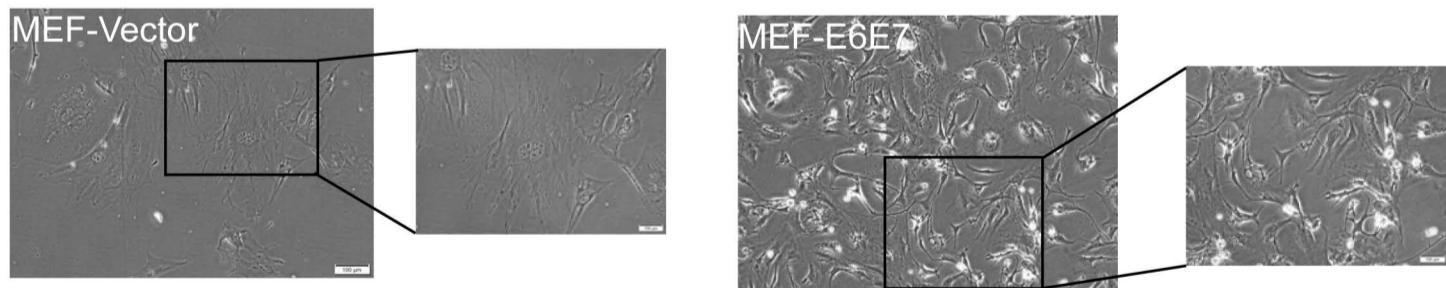
S1C



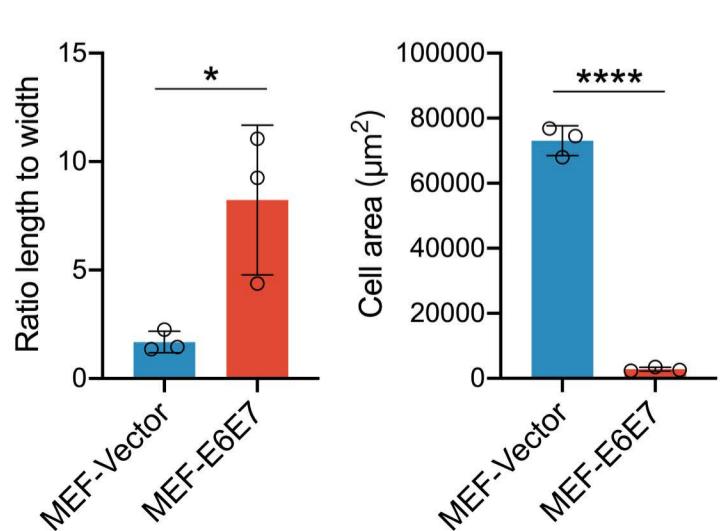
S1D



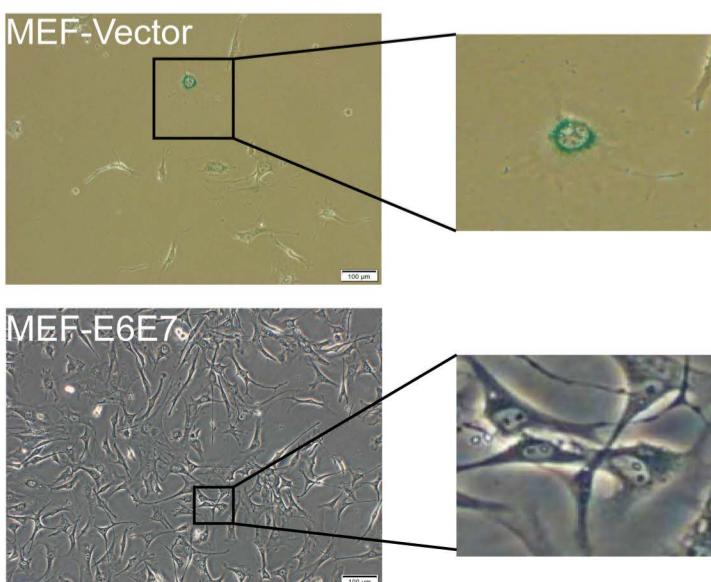
S1E



S1F

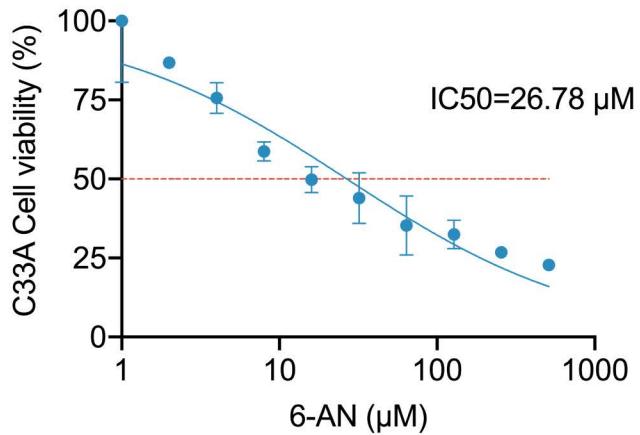


S1G

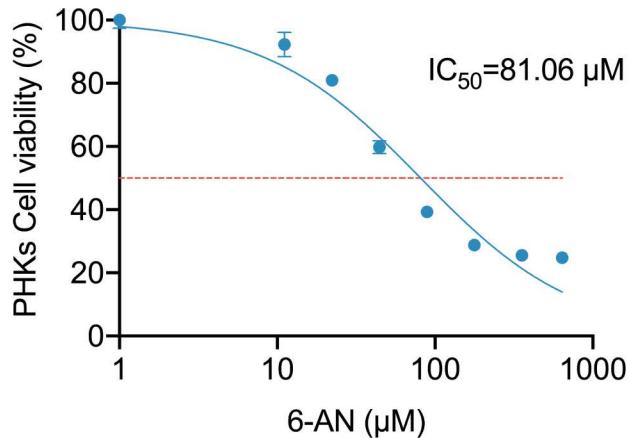


Supplementary Figure 2

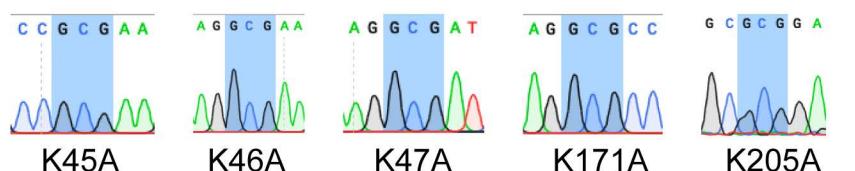
S2A



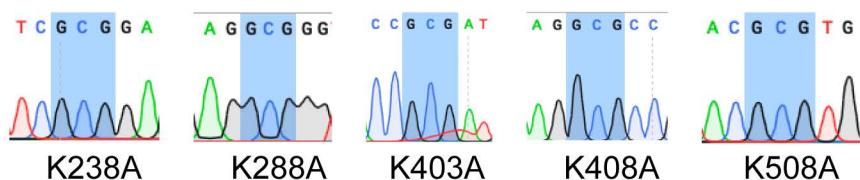
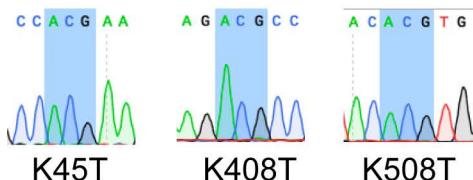
S2B



S2C

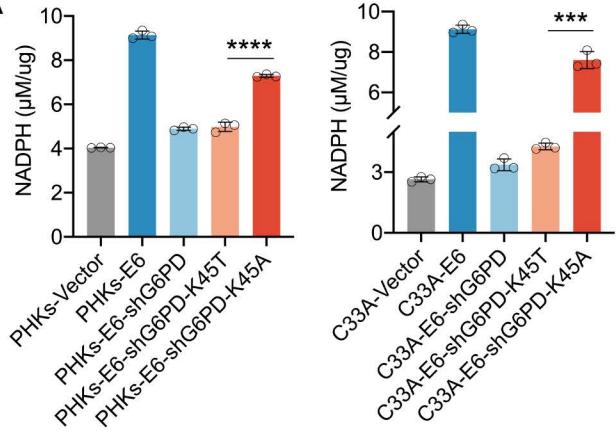


S2D

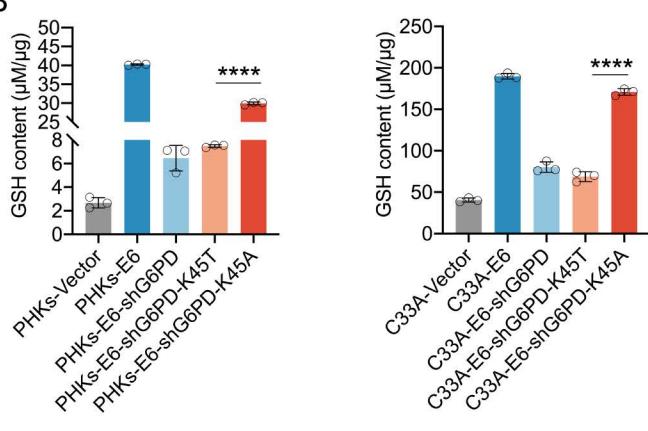


Supplementary Figure 3

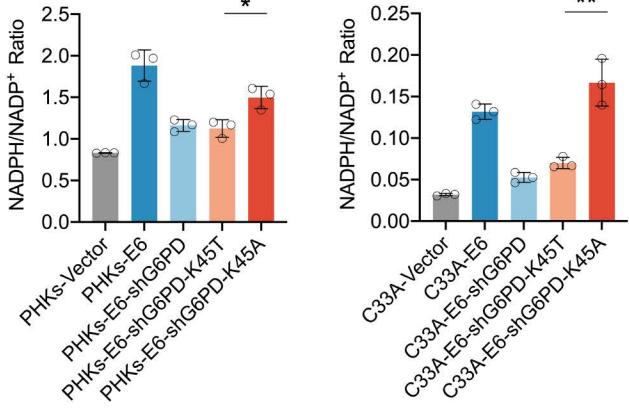
S3A



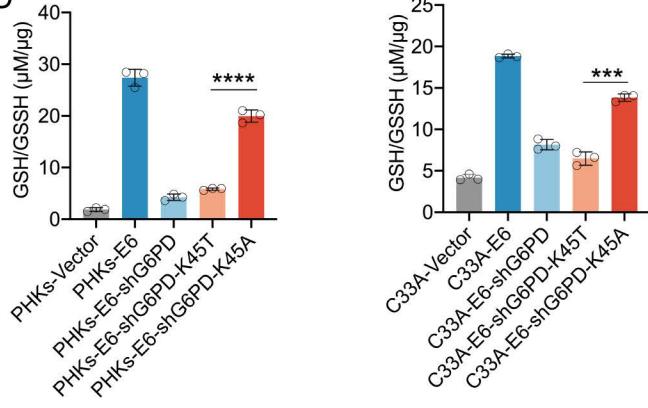
S3B



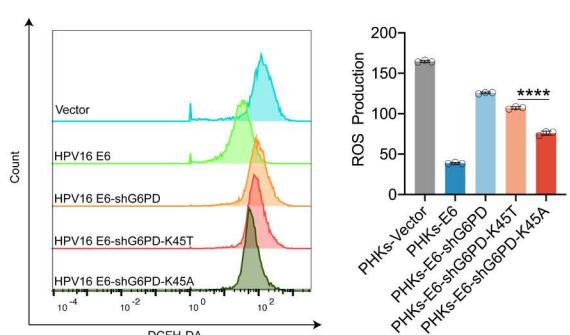
S3C



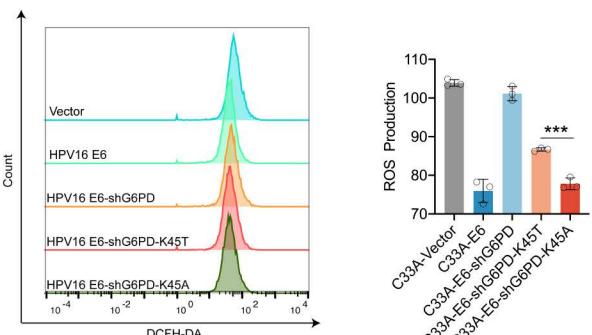
S3D



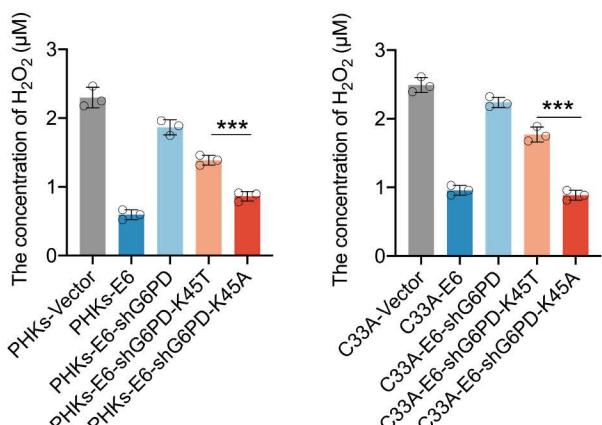
S3E



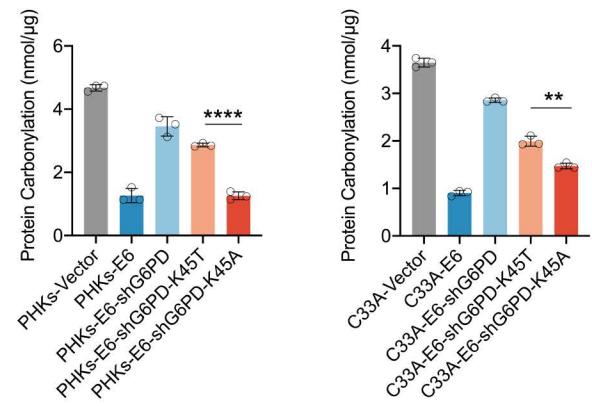
S3F



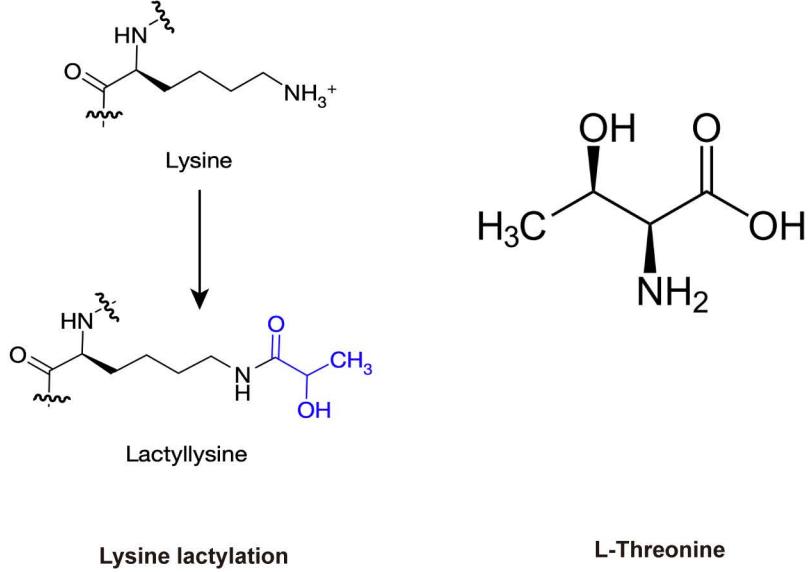
S3G



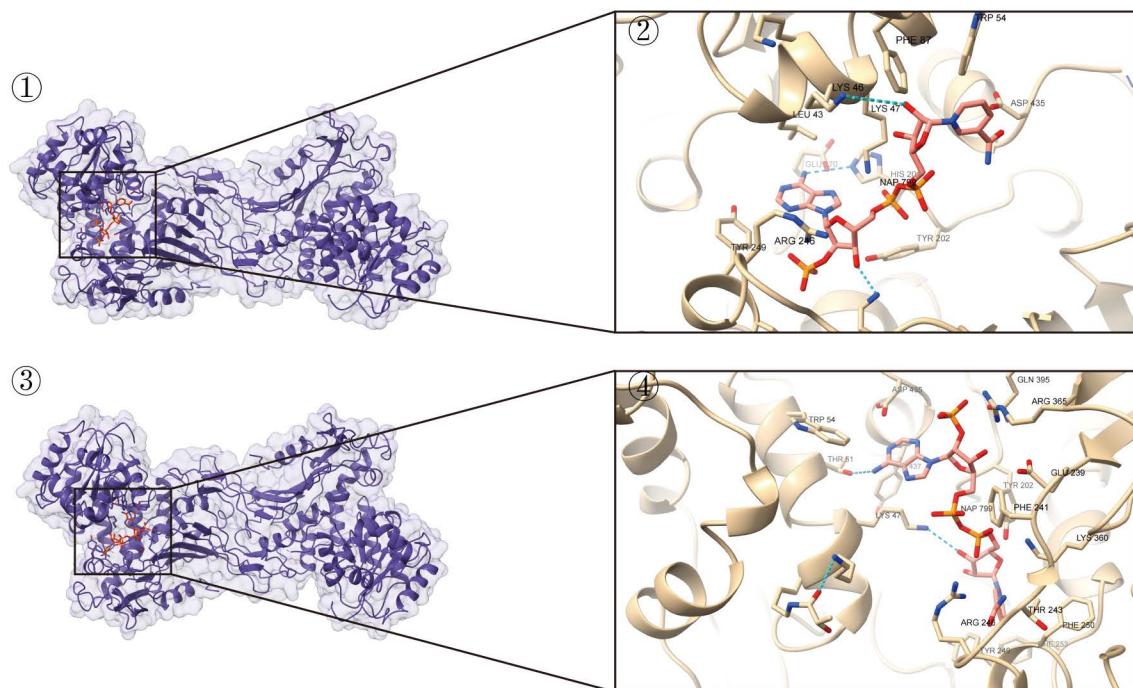
S3H



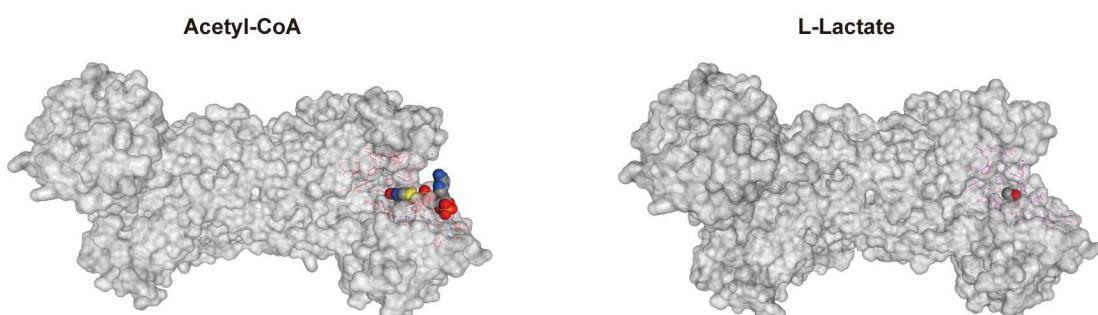
S4A



S4B



S4C



Ligand	Receptor	Docking Score
Acetyl-CoA	G6PD	175.97
L-Lactate	G6PD	208.85

Table represents the L-Lactate/L-Lactate—DCBLD1 interacting atoms and dock score analyzed by online tool CB-Dock2 (<https://cadd.labshare.cn/cb-dock2/index.php>).