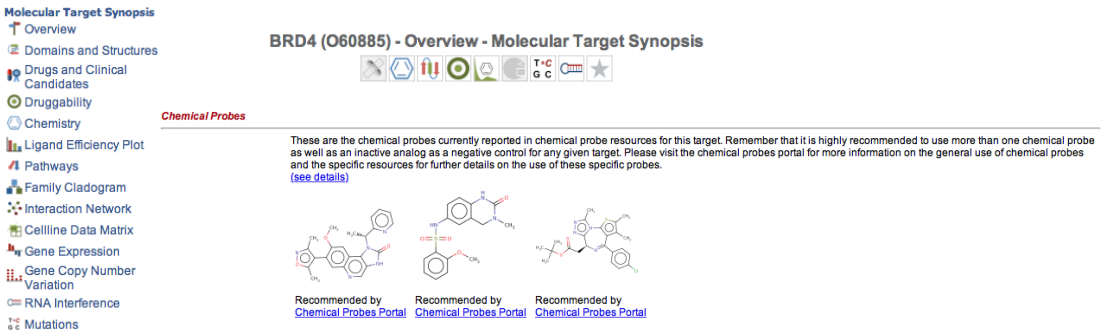
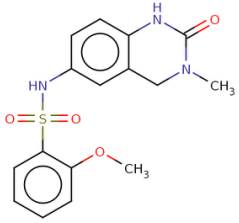
Selecting the best chemical probe for a target



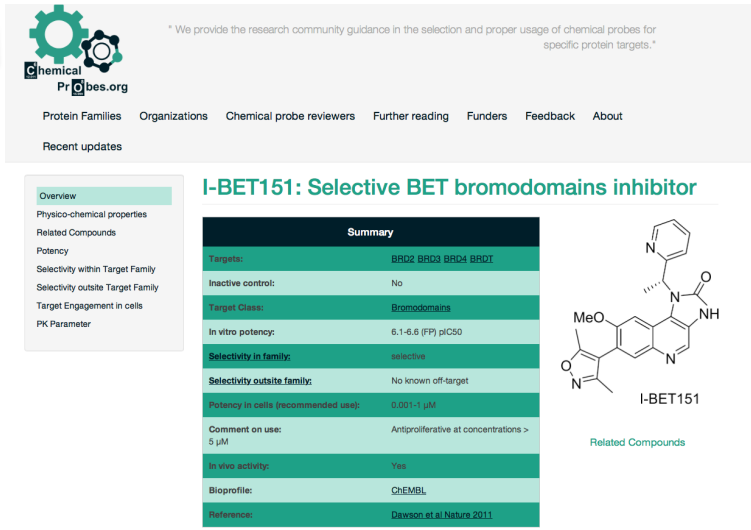
On the molecular target synposis page, scroll down to ‘Chemical Probes’. This shows chemical probes currently reported in chemical probe resources for this target. Remember that it is highly recommended to use more than one chemical probe as well as an inactive analog as a negative control for any given target.



1016603 - Overview - Compound Synopsis

Synonyms	
0NS	
Molecular Formula	
C16H17N3O4S	
IUPAC	
'2-methoxy-N-(3-methyl-2-oxo-1,2,3,4-tetrahydroquinazolin-6-yl)benzene-1-sulfona...	
Standard InChI	
InChI=1S/C16H17N3O4S/c1-19-10-11-9-12(7-8-13(11)17-16(19)20)18-24(21,22)15-6-4-3...	
SD File	
SciTegic10121214212D 24 26 0 0 0 0 999 V2000 -7.5533 0...	
Smiles	
COc1ccccc1S(=O)(=O)Nc2ccc3NC(=O)N(C)Cc3c2	
Cross References	
PDB Het Code	0NS
ChEMBL	CHEMBL2179387
Targets Recommended as a Probe	
O60885, Q15059, P25440, Q58F21	

Clicking on a chemical probe structure takes you to its compound synopsis page where you can explore other compounds with similar scaffolds, its protein affinity profile and more.



Clicking the link to the Chemical Probes Portal (http://www.chemicalprobes.org) will take you to the Chemical Probes Portal entry for this probe.