Supporting Information: Elemental Analysis of Target Compounds

Structure-based Design of Novel Chk1 Inhibitors: Insights into Hydrogen Bonding and Protein-Ligand Affinity

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Compound	Formulae	Elemental Analysis					
		Calculated			Found		
		C	H	N	C	H	N
2	C20 H17 N3 O2	72.68	5.18	12.63	72.49	5.17	12.67
3	C21 H19 N3 O2	71.73	5.63	11.88	73.03	5.54	12.16
4	C20 H15 N3 O3	72.68	5.18	12.63	72.49	5.17	12.67
5	C20 H18 N4 O	72.09	5.49	16.90	72.71	5.49	16.95
6	C21 H20 N4 O	72.55	5.83	16.20	73.23	5.85	16.26
7	C21 H20 N4 O2	67.94	5.46	15.10	69.98	5.59	15.54