

Characterization data of 6a-r

N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)benzamide (6a): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr) 3355 (Amide NH stretch), 3055 (Aromatic -CH stretch), 1665 (Amide -C=O stretch), 1653, 1600 (Imine -C=N stretch), 1556 (-CH bending), 1541, 1484, 699; ¹H NMR (500 MHz, DMSO) δ 10.58 (s, 1H, D₂O exchangeable, Amide NH), 8.90 (d, *J* = 6.7 Hz, 1H, ArH), 8.77 (s, 1H, ArH), 8.56 (s, 1H, ArH), 8.03 (d, *J* = 7.2 Hz, 2H, ArH), 7.97 – 7.89 (m, 2H, ArH), 7.79 (t, *J* = 9.6 Hz, 2H, ArH), 7.63 (dd, *J* = 13.3, 6.0 Hz, 1H, ArH), 7.60 – 7.55 (m, 3H, ArH), 7.48 (t, *J* = 6.6 Hz, 1H, ArH); ¹³C NMR (500 MHz, DMSO) δ 166.00 (Amide, -C=O), 145.27 (ArC=N), 144.78 (ArC), 140.10 (ArC), 135.37 (ArC), 134.80 (ArC), 132.06 (ArC), 129.42 (ArC), 128.86 (ArC), 128.15 (ArC), 127.41 (ArC), 125.45 (ArC), 121.43 (ArC), 120.15 (ArC), 118.14 (ArC), 117.09 (ArC), 112.73 (ArC), 109.60 (ArC); LC/MS Calculated C₂₀H₁₅N₃O [M+1]⁺: m/z 313.36; observed, 314.00.

4-Fluoro-N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)benzamide (6b): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr) 3351 (Amide NH stretch), 3069 (Aromatic -CH stretch), 1658 (Amide -C=O stretch), 1604 (Imine -C=N stretch), 1559 (-CH bending), 755 (-CF stretch); ¹H NMR (500 MHz, DMSO) δ 10.63 (s, 1H, Amide NH), 8.89 (d, *J* = 6.7 Hz, 1H, ArH), 8.75 (s, 1H, ArH), 8.54 (t, *J* = 1.7 Hz, 1H, ArH), 8.13 (dd, *J* = 8.8, 5.5 Hz, 2H, ArH), 7.95 (d, *J* = 9.0 Hz, 1H, ArH), 7.91 – 7.85 (m, 1H, ArH), 7.83 – 7.77 (m, 2H, ArH), 7.58 (t, *J* = 7.9 Hz, 1H, ArH), 7.46 – 7.37 (m, 3H, ArH); ¹³C NMR (500 MHz, DMSO) δ 166.82 (-CF), 165.09 (Amide, -C=O), 140.37 (ArC=N), 132.61 (ArC), 132.53 (ArC), 131.52 (ArC), 131.03 (ArC), 130.96 (ArC), 130.09 (ArC), 128.88 (ArC), 121.97 (ArC), 118.45 (ArC), 116.18 (ArC), 115.96 (ArC), 114.06 (ArC), 111.03 (ArC); LC/MS Calculated C₂₀H₁₄FN₃O [M+1]⁺: m/z 331.35, observed, 332.10.

4-Chloro-N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)benzamide (6c): Nature and colour: White crystalline solid; FT-IR (cm⁻¹) (KBr); 3353 (Amide NH stretch), 3069 (Aromatic -CH

stretch), 1661 (Amide -C=O stretch), 1599 (Imine -C=N stretch), 1550 (-CH bending), 750 (-CCl stretch); ¹H NMR (500 MHz, DMSO) δ 10.42 (s, 1H, Amide NH), 8.56 (dt, *J* = 6.8, 1.1 Hz, 1H, ArH), 8.47 (s, 1H, ArH), 8.38 (s, 1H, ArH), 8.05 (d, *J* = 8.6 Hz, 2H, ArH), 7.77 (dd, *J* = 8.5, 1.6 Hz, 1H, ArH), 7.72 – 7.68 (m, 1H, ArH), 7.64 (d, *J* = 8.6 Hz, 2H, ArH), 7.59 (d, *J* = 9.1 Hz, 1H, ArH), 7.43 (t, *J* = 7.9 Hz, 1H, ArH), 7.26 (ddd, *J* = 9.0, 6.7, 1.2 Hz, 1H, ArH), 6.91 (td, *J* = 6.7, 1.1 Hz, 1H, ArH); ¹³C NMR (500 MHz, DMSO) δ 164.89 (Amide, -C=O), 145.28 (ArC=N), 144.73, 139.91 (-CCl), 136.92 (ArC), 134.85 (ArC), 134.06 (ArC), 130.13 (ArC), 129.47 (ArC), 128.95 (ArC), 127.42 (ArC), 125.48 (ArC), 121.59 (ArC), 120.18 (ArC), 118.18 (ArC), 117.10 (ArC), 112.75 (ArC), 109.63 (ArC); LC/MS Calculated C₂₀H₁₄ClN₃O [M+1]⁺: m/z 347.80, observed, 348.00.

4-Bromo-N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)benzamide (6d): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr) 3376 (Amide NH stretch), 3134 (Aromatic -CH stretch), 1656 (Amide -C=O stretch), 1614 (Imine -C=N stretch), 1558 (-CH bending), 739 (-CBr stretch); ¹H NMR (500 MHz, DMSO) δ 10.60 (s, 1H, Amide NH), 8.81 (d, *J* = 6.7 Hz, 1H, ArH), 8.67 (s, 1H, ArH), 8.52 (t, *J* = 1.8 Hz, 1H, ArH), 8.01 – 7.94 (m, 2H, ArH), 7.86 (d, *J* = 9.0 Hz, 1H, ArH), 7.82 – 7.71 (m, 5H, ArH), 7.55 (t, *J* = 7.9 Hz, 1H, ArH), 7.33 (t, *J* = 6.5 Hz, 1H, ArH); ¹³C NMR (500 MHz, DMSO) δ 165.12 (Amide, -C=O), 143.89 (ArC=N), 140.06 (ArC), 134.29 (ArC), 131.91 (ArC), 130.34 (ArC), 129.74 (ArC), 128.08 (ArC), 125.95 (-CBr), 121.81 (ArC), 118.32 (ArC), 115.71 (ArC), 114.28 (ArC), 110.29 (ArC); LC/MS Calculated C₂₀H₁₄BrN₃O [M+1]⁺: m/z 392.26, observed, 393.90;

N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)-4-(trifluoromethyl)benzamide (6e): Nature and colour: White crystalline solid; FT-IR (cm⁻¹) (KBr); 3269 (Amide NH stretch), 3061 (Aromatic -CH stretch), 1666 (Amide -C=O stretch), 1615 (Imine -C=N stretch), 1557 (-CH bending), 750; ¹H NMR (500 MHz, DMSO) δ 10.57 (s, 1H, Amide NH), 8.57 (d, *J* = 6.8 Hz, 1H, ArH), 8.48 (s, 1H, ArH), 8.40 (s, 1H, ArH), 8.21 (d, *J* = 8.1 Hz, 2H, ArH), 7.94 (d, *J* = 8.2 Hz, 2H,

ArH), 7.77 (dd, $J = 8.1, 1.2$ Hz, 1H, ArH), 7.71 (d, $J = 7.9$ Hz, 1H, ArH), 7.59 (d, $J = 9.1$ Hz, 1H, ArH), 7.45 (t, $J = 7.9$ Hz, 1H, ArH), 7.27 (ddd, $J = 9.0, 6.7, 1.1$ Hz, 1H, ArH), 6.92 (td, $J = 6.8, 1.0$ Hz, 1H, ArH); ^{13}C NMR (500 MHz, DMSO) δ 164.86 (Amide, -C=O), 145.23 (ArC=N), 144.55 (ArC), 139.77 (ArC), 134.79 (ArC), 131.99 (ArC), 129.54 (ArC), 127.45 (ArC), 125.89 (-CF₃), 121.79 (ArC), 120.26 (ArC), 118.22 (ArC), 117.04 (ArC), 112.83 (ArC), 109.69 (ArC); LC/MS Calculated C₂₁H₁₄F₃N₃O [M+1]⁺: m/z 381.36, observed, 382.10.

N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)-4-methylbenzamide (6f): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr); 3268 (Amide NH stretch), 3066 (Aromatic -CH stretch), 2985 (Aliphatic CH stretch), 1669 (Amide -C=O stretch), 1618 (Imine -C=N stretch), 1559 (-CH bending), 756; ^1H NMR (500 MHz, DMSO) δ 10.26 (s, 1H, Amide NH), 8.56 (d, $J = 8.8$ Hz, 1H, ArH), 8.48 (s, 1H, ArH), 8.38 (s, 1H, ArH), 7.94 (d, $J = 8.2$ Hz, 2H, ArH), 7.80 – 7.76 (m, 1H, ArH), 7.70 – 7.67 (m, 1H, ArH), 7.59 (d, $J = 9.1$ Hz, 1H, ArH), 7.42 (t, $J = 7.9$ Hz, 1H, ArH), 7.36 (d, $J = 7.9$ Hz, 2H, ArH), 7.26 (ddd, $J = 9.0, 6.7, 1.2$ Hz, 1H, ArH), 6.91 (t, $J = 7.3$ Hz, 1H, ArH), 2.41 (s, 3H, CH₃); ^{13}C NMR (500 MHz, DMSO) δ 166.39 (Amide, -C=O), 140.91 (ArC=N), 138.23 (ArCCH₃), 135.01 (ArC), 132.85 (ArC), 130.23 (ArC), 129.51 (ArC), 128.83 (ArC), 125.43 (ArC), 122.84 (ArC), 118.53 (ArC), 117.55 (ArC), 112.68 (ArC), 111.52 (ArC), 21.46 (CH₃); LC/MS Calculated C₂₁H₁₇N₃O [M+1]⁺: m/z 327.39, observed, 328.10.

N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)-4-methoxybenzamide (6g): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr); 3270 (Amide NH stretch), 3069 (Aromatic -CH stretch), 2965 (Aliphatic CH stretch), 1664 (Amide -C=O stretch), 1619 (Imine -C=N stretch), 1600 (-CH bending), 759; ^1H NMR (500 MHz, DMSO) δ 10.44 (s, 1H, Amide NH), 8.92 (d, $J = 6.7$ Hz, 1H, ArH), 8.78 (s, 1H, ArH), 8.56 (s, 1H, ArH), 8.05 (d, $J = 8.9$ Hz, 2H, ArH), 7.94 (dt, $J = 19.1, 6.1$ Hz, 2H, ArH), 7.82 – 7.76 (m, 2H, ArH), 7.58 (t, $J = 7.9$ Hz, 1H, ArH), 7.50 (t, $J = 7.2$ Hz, 1H, ArH), 7.10 (d, $J = 8.9$ Hz, 2H, ArH), 3.86 (s, 3H, OCH₃); ^{13}C

NMR (500 MHz, DMSO) δ 165.63 (Amide, -C=O), 162.25 (ArCOCH₃), 140.80 (ArC=N), 136.42 (ArC), 133.47 (ArC), 130.22 (ArC), 129.48 (ArC), 127.01 (ArC), 122.76 (ArC), 121.85 (ArC), 118.49 (ArC), 117.48 (ArC), 114.16 (ArC), 112.81 (ArC), 111.56 (ArC), 55.96 (OCH₃); LC/MS Calculated C₂₁H₁₇N₃O₂ [M+1]⁺: m/z 343.39, observed 344.10.

N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)-4-nitrobenzamide (6h): Nature and colour: Bright yellow crystalline solid; FT-IR (cm⁻¹) (KBr); 3313 (Amide NH stretch), 3076 (Aromatic -CH stretch), 1645 (Amide -C=O stretch), 1602 (Imine -C=N stretch), 1515 (Nitro -NO₂ symmetric stretch), 1346 (Nitro -NO₂ asymmetric stretch), 764; ¹H NMR (500 MHz, DMSO) δ 10.67 (s, 1H, Amide NH), 8.56 (d, *J* = 6.8 Hz, 1H, ArH), 8.48 (s, 1H, ArH), 8.42 – 8.37 (m, 3H, ArH), 8.28 – 8.22 (m, 2H, ArH), 7.78 (dd, *J* = 8.0, 1.2 Hz, 1H, ArH), 7.74 – 7.70 (m, 1H, ArH), 7.59 (d, *J* = 9.6 Hz, 1H, ArH), 7.45 (t, *J* = 7.9 Hz, 1H, ArH), 7.27 (ddd, *J* = 9.0, 6.7, 1.2 Hz, 1H, ArH), 6.92 (td, *J* = 6.7, 1.1 Hz, 1H, ArH); ¹³C NMR (500 MHz, DMSO) δ 164.26 (Amide, -C=O), 149.66 (ArCNO₂), 145.30 (ArC=N), 144.50 (ArC), 141.02 (ArC), 139.66 (ArC), 136.86 (ArC), 134.85 (ArC), 131.17 (ArC), 129.72 (ArC), 127.45 (ArC), 125.60 (ArC), 124.21 (ArC), 121.92 (ArC), 120.24 (ArC), 118.23 (ArC), 117.06 (ArC), 112.82 (ArC), 109.70 (ArC); LC/MS Calculated C₂₀H₁₄N₄O₃ [M+1]⁺: m/z 358.36, observed, 359.10.

3-Fluoro-N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)benzamide (6i): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr); 3372 (Amide NH stretch), 3074 (Aromatic -CH stretch), 1667 (Amide -C=O stretch), 1618 (Imine -C=N stretch), 1591 (-CH bending), 750 (-CF stretch); ¹H NMR (500 MHz, DMSO) δ 10.64 (s, 1H, Amide NH), 8.88 (d, *J* = 6.7 Hz, 1H, ArH), 8.75 (s, 1H, ArH), 8.54 (s, 1H, ArH), 7.96 – 7.83 (m, 4H, ArH), 7.78 (dd, *J* = 9.2, 8.1 Hz, 2H, ArH), 7.62 (dt, *J* = 21.6, 8.0 Hz, 2H, ArH), 7.52 – 7.42 (m, 2H, ArH); ¹³C NMR (500 MHz, DMSO) δ 163.38 (Amide, -C=O), 161.44 (-CF), 140.90 (ArC=N), 140.31 (ArC), 137.28 (ArC), 136.17 (ArC), 133.50 (ArC), 131.16 (ArC), 130.26 (ArC), 129.49 (ArC), 127.63 (ArC),

124.53 (ArC), 122.94 (ArC), 118.65 (ArC), 117.94 (ArC), 115.20 (ArC), 112.74 (ArC), 111.59 (ArC); LC/MS Calculated C₂₀H₁₄FN₃O [M+1]⁺: m/z 331.35, observed 332.10.

3-Chloro-N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)benzamide (6j): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr); 3362 (Amide NH stretch), 3075 (Aromatic -CH stretch), 1666 (Amide -C=O stretch), 1617 (Imine -C=N stretch), 1596 (-CH bending), 746 (-CCl stretch); ¹H NMR (500 MHz, DMSO) δ 10.69 (s, 1H, Amide NH), 8.90 (d, *J* = 6.7 Hz, 1H, ArH), 8.77 (s, 1H, ArH), 8.53 (s, 1H, ArH), 8.08 (t, *J* = 1.8 Hz, 1H, ArH), 8.02 – 7.98 (m, 1H, ArH), 7.95 (d, *J* = 8.9 Hz, 1H, ArH), 7.93 – 7.88 (m, 1H, ArH), 7.79 (dd, *J* = 7.9 Hz, 2H, ArH), 7.71 (ddd, *J* = 8.0 Hz, 1H, ArH), 7.61 (td, *J* = 7.9, 4.3 Hz, 2H, ArH), 7.47 (t, *J* = 6.6 Hz, 1H, ArH); ¹³C NMR (500 MHz, DMSO) δ 164.80 (Amide, -C=O), 140.96 (ArC=N), 140.30 (ArC), 136.94 (ArC), 133.74 (ArCCl), 132.11 (ArC), 130.95 (ArC), 129.50 (ArC), 128.01 (ArC), 127.11 (ArC), 122.94 (ArC), 122.37 (ArC), 118.64 (ArC), 117.52 (ArC), 112.74 (ArC), 111.61 (ArC); LC/MS Calculated C₂₀H₁₄ClN₃O [M+1]⁺: m/z 347.80; observed, 346.99 (deconvoluted).

3-Bromo-N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)benzamide (6k): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr) 3258 (Amide NH stretch), 3125 (Heteroaromatic -CH stretch), 3038 (Aromatic -CH stretch), 1640 (Amide -C=O stretch), 1613 (Imine -C=N stretch), 1554 (-CH bending); ¹H NMR (500 MHz, DMSO) δ 10.49 (s, 1H, Amide NH), 8.62 (d, *J* = 6.7 Hz, 1H, ArH), 8.46 (d, *J* = 10.2 Hz, 2H, ArH), 8.22 (t, *J* = 1.8 Hz, 1H, ArH), 8.05 – 7.98 (m, 1H, ArH), 7.82 (ddd, *J* = 8.0, 1.8, 0.8 Hz, 1H, ArH), 7.78 (dd, *J* = 8.1, 1.2 Hz, 1H, ArH), 7.71 (d, *J* = 7.8 Hz, 1H, ArH), 7.65 (d, *J* = 9.0 Hz, 1H, ArH), 7.53 (t, *J* = 7.9 Hz, 1H, ArH), 7.46 (t, *J* = 7.9 Hz, 1H, ArH), 7.41 – 7.36 (m, 1H, ArH), 7.02 (t, *J* = 6.7 Hz, 1H, ArH); ¹³C NMR (500 MHz, DMSO) δ 164.47 (Amide, -C=O), 144.52 (ArC=N), 139.92 (ArC), 137.84 (ArC), 134.84 (ArC), 131.15 (-CBr), 130.77 (ArC), 129.63 (ArC), 127.79

(ArC), 122.19 (ArC), 121.78 (ArC), 120.65 (ArC), 118.26 (ArC), 116.33 (ArC), 113.59 (ArC), 110.01 (ArC); LC/MS Calculated C₂₀H₁₄BrN₃O [M+1]⁺: m/z 392.26, observed, 392.00.

N-(3-(imidazo[1,2-a]pyridin-2-yl)phenyl)-3-(trifluoromethyl)benzamide (6l): Nature and colour: White crystalline solid; FT-IR (cm⁻¹) (KBr); 3267 (Amide NH stretch), 3064 (Aromatic -CH stretch), 1667 (Amide -C=O stretch), 1614 (Imine -C=N stretch), 1537 (-CH bending), 773 (-CF stretch); ¹H NMR (500 MHz, DMSO) δ 10.57 (s, 1H, Amide NH), 8.56 (d, *J* = 7.9 Hz, 1H, ArH), 8.46 (s, 1H, ArH), 8.38 (d, *J* = 10.2 Hz, 2H, ArH), 8.32 (d, *J* = 7.8 Hz, 1H, ArH), 7.99 (d, *J* = 7.8 Hz, 1H, ArH), 7.81 (t, *J* = 7.8 Hz, 2H, ArH), 7.71 (d, *J* = 7.9 Hz, 1H, ArH), 7.58 (d, *J* = 9.7 Hz, 1H, ArH), 7.45 (t, *J* = 7.9 Hz, 1H, ArH), 7.27 (dd, *J* = 6.7, 1.2 Hz, 1H, ArH), 6.91 (td, *J* = 6.7, 1.1 Hz, 1H, ArH); ¹³C NMR (500 MHz, DMSO) δ 164.47 (Amide, -C=O), 145.29 (ArC=N), 144.66 (ArC), 139.76 (ArC), 136.19 (ArC), 134.89 (ArC), 132.38 (ArC), 130.22 (ArCCF₃), 129.79 (ArC), 127.43 (ArC), 125.57 (ArC), 124.73 (ArC), 123.40 (-CF), 121.75 (ArC), 120.24 (ArC), 118.28 (ArC), 117.09 (ArC), 112.77 (ArC), 109.66 (ArC); LC/MS Calculated C₂₀H₁₄ClN₃O [M+1]⁺: m/z 381.36; observed 381.09 (deconvoluted).

N-(3-(imidazo[1,2-a]pyridin-2-yl)phenyl)-3-methylbenzamide (6m): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr); 3375 (Amide NH stretch), 3050 (Aromatic -CH stretch), 2922 (Aliphatic CH stretch), 1667 (Amide -C=O stretch), 1618 (Imine -C=N stretch), 1591 (-CH bending), 742; ¹H NMR (500 MHz, DMSO) δ 10.55 (s, 1H, Amide NH), 8.92 (d, *J* = 6.7 Hz, 1H, ArH), 8.78 (s, 1H, ArH), 8.54 (d, *J* = 1.6 Hz, 1H, ArH), 7.99 – 7.90 (m, 2H, ArH), 7.85 (s, 1H, ArH), 7.84 – 7.77 (m, 3H, ArH), 7.59 (t, *J* = 7.9 Hz, 1H, ArH), 7.49 (dd, *J* = 11.6, 5.7 Hz, 1H, ArH), 7.45 (dd, *J* = 4.5, 2.0 Hz, 2H, ArH), 2.44 – 2.41 (m, 3H, ArH); ¹³C NMR (500 MHz, DMSO) δ 166.39 (Amide, -C=O), 140.91 (ArC=N), 138.23 (ArCCH₃), 135.01 (ArC), 132.85 (ArC), 130.23 (ArC), 129.51 (ArC), 128.83 (ArC), 125.43 (ArC), 122.84 (ArC), 118.53 (ArC), 117.55 (ArC), 112.68 (ArC), 111.52 (ArC), 21.46 (CH₃); LC/MS Calculated C₂₁H₁₇N₃O [M+1]⁺: m/z 327.39, observed, 328.10.

N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)-3-methoxybenzamide (6n): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr); 3319 (Amide NH stretch), 3074 (Aromatic -CH stretch), 2967 (Aliphatic CH stretch), 1666 (Amide -C=O stretch), 1617 (Imine -C=N stretch), 1586 (-CH bending), 743; ¹H NMR (500 MHz, DMSO) δ 10.59 (s, 1H, Amide NH), 8.92 (d, *J* = 6.7 Hz, 1H, ArH), 8.78 (s, 1H, ArH), 8.55 (s, 1H, ArH), 7.97 (d, *J* = 8.9 Hz, 1H, ArH), 7.95 – 7.90 (m, 1H, ArH), 7.81 (t, *J* = 6.9 Hz, 2H, ArH), 7.65 – 7.60 (m, 1H, ArH), 7.58 (dd, *J* = 5.0, 2.7 Hz, 2H, ArH), 7.52 – 7.45 (m, 2H, ArH), 7.20 (dd, *J* = 8.1, 2.3 Hz, 1H, ArH), 3.87 (s, 3H, OCH₃); ¹³C NMR (500 MHz, DMSO) δ 165.98 (Amide, -C=O), 159.68 (ArCOCH₃), 140.98 (ArC=N), 140.53 (ArC), 136.36 (ArC), 133.55 (ArC), 130.22 (ArC), 129.50 (ArC), 122.96 (ArC), 122.15 (ArC), 120.49 (ArC), 118.66 (ArC), 117.53 (ArC), 113.51 (ArC), 112.71 (ArC), 111.58 (ArC), 55.89 (OCH₃); LC/MS Calculated C₂₁H₁₇N₃O₂ [M+1]⁺: m/z 343.39, observed, 344.10.

N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)-3-nitrobenzamide (6o): Light yellow crystalline solid; FT-IR (cm⁻¹) (KBr); 3241 (Amide NH stretch), 3084 (Aromatic -CH stretch), 1668 (Amide -C=O stretch), 1615 (Imine -C=N stretch), 1552 (Nitro –NO₂ symmetric stretch), 1346 (Nitro –NO₂ asymmetric stretch), 752; ¹H NMR (500 MHz, DMSO) δ 10.73 (s, 1H, Amide NH), 8.87 (s, 1H, ArH), 8.61 (d, *J* = 6.8 Hz, 2H, ArH), 8.49-8.44 (m, 2H, ArH), 8.36 (d, *J* = 10.3 Hz, 1H, ArH), 7.87 (t, *J* = 8.0 Hz, 1H, ArH), 7.82 (t, *J* = 8.0 Hz, 1H, ArH), 7.73 (d, *J* = 7.8 Hz, 1H, ArH), 7.64 (d, *J* = 9.0 Hz, 1H, ArH), 7.48 (t, *J* = 7.9 Hz, 1H, ArH), 7.39 – 7.33 (m, 1H, ArH), 7.00 (t, *J* = 6.6 Hz, 1H, ArH); ¹³C NMR (500 MHz, DMSO) δ 163.83 (Amide, -C=O), 148.25 (ArCNO₂), 144.68 (ArC=N), 139.73 (ArC), 134.70 (ArC), 131.01 (ArC), 130.68 (ArC), 129.66 (ArC), 127.81 (ArC), 122.94 (ArC), 121.95 (ArC), 120.65 (ArC), 113.45 (ArC), 109.98 (ArC); LC/MS Calculated C₂₀H₁₄N₄O₃ [M+1]⁺: m/z 358.36, observed, 359.00.

2-Fluoro-N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)benzamide (6p): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr); 3287 (Amide NH stretch), 3072 (Aromatic -CH stretch), 1656 (Amide -C=O stretch), 1612 (Imine -C=N stretch), 1591 (-CH bending), 745 (-CF stretch); ¹H NMR (500 MHz, DMSO) δ 10.70 (s, 1H, Amide NH), 8.87 (d, *J* = 6.7 Hz, 1H, ArH), 8.75 (s, 1H, ArH), 8.49 (s, 1H, ArH), 7.93 (d, *J* = 8.9 Hz, 1H, ArH), 7.86 (t, *J* = 7.8 Hz, 1H, ArH), 7.78 (d, *J* = 7.8 Hz, 1H, ArH), 7.75-7.67 (m, 2H, ArH), 7.65 -7.55 (m, 2H, ArH), 7.40 (ddd, *J* = 16.0, 10.9, 3.8 Hz, 3H, ArH); ¹³C NMR (500 MHz, DMSO) δ 164.89 (-CF), 162.34 (Amide, -C=O), 145.28 (ArC=N), 144.75 (ArC), 140.01 (ArC), 134.82 (ArC), 131.80 (ArC), 130.94 (ArC), 129.45 (ArC), 127.41 (ArC), 125.48 (ArC), 121.50 (ArC), 120.18 (ArC), 118.17 (ArC), 117.09 (ArC), 115.90 (ArC), 112.74 (ArC), 109.62 (ArC); LC/MS Calculated C₂₀H₁₄FN₃O [M+1]⁺: m/z 331.35; observed 330.89 (deconvoluted).

N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)-1-naphthamide (6q): Nature and colour: Off-white crystalline solid; FT-IR (cm⁻¹) (KBr); 3239 (Amide NH stretch), 3063 (Aromatic -CH stretch), 1651 (Amide -C=O stretch), 1615 (Imine -C=N stretch), 1593 (-CH bending), 740; ¹H NMR (500 MHz, DMSO) δ 10.67 (s, 1H, Amide NH), 8.57 (d, *J* = 7.0 Hz, 2H, ArH), 8.40 (s, 1H, ArH), 8.28 – 8.23 (m, 1H, ArH), 8.10 (d, *J* = 8.3 Hz, 1H, ArH), 8.05 (dd, *J* = 6.7, 2.7 Hz, 1H, ArH), 7.81 (dd, *J* = 6.9, 0.7 Hz, 1H, ArH), 7.73 (dd, *J* = 7.7, 1.2 Hz, 2H, ArH), 7.67 – 7.58 (m, 4H, ArH), 7.45 (t, *J* = 7.9 Hz, 1H, ArH), 7.26 (ddd, *J* = 9.1, 6.7, 1.2 Hz, 1H, ArH), 6.91 (td, *J* = 6.7, 1.1 Hz, 1H, ArH); ¹³C NMR (500 MHz, DMSO) δ 167.85 (Amide, -C=O), 145.29 (ArC=N), 144.80 (ArC), 140.24 (ArC), 133.66 (ArC), 130.61 (ArC), 129.55 (ArC), 128.82 (ArC), 127.49 (ArC), 127.42 (ArC), 126.84 (ArC), 125.99 (ArC), 125.64 (ArC), 125.54 (ArC), 125.46 (ArC), 117.13 (ArC), 112.74 (ArC), 109.66 (ArC); LC/MS Calculated C₂₄H₁₇N₃O [M+1]⁺: m/z 363.42; observed, 364.10.

N-(3-(imidazo[1,2-*a*]pyridin-2-yl)phenyl)furan-2-carboxamide (6r): Nature and colour: Dark brown crystalline solid; FT-IR (cm⁻¹) (KBr); 3283 (Amide NH stretch), 3127

(heteroaromatic -CH stretch), 1650 (Amide -C=O stretch), 1613 (Imine -C=N stretch), 1579 (-CH bending), 751; ¹H NMR (500 MHz, DMSO) δ 10.26 (s, 1H, Amide NH), 8.56 (d, *J* = 8.9 Hz, 1H, ArH), 8.42 (s, 1H, ArH), 8.37 (s, 1H, ArH), 7.96 (dd, *J* = 1.6, 0.7 Hz, 1H, ArH), 7.75 (ddd, *J* = 8.1, 2.0, 0.8 Hz, 1H, ArH), 7.71-7.66 (m, 1H, ArH), 7.59 (d, *J* = 9.1 Hz, 1H, ArH), 7.41 (dd, *J* = 10.1, 5.7 Hz, 2H, ArH), 7.26 (ddd, *J* = 9.0, 6.7, 1.2 Hz, 1H, ArH), 6.91 (td, *J* = 6.7, 1.1 Hz, 1H, ArH), 6.73 (dd, *J* = 3.5, 1.7 Hz, 1H, ArH); ¹³C NMR (500 MHz, DMSO) δ 156.76 (Amide, -C=O), 147.96 (ArC=N), 146.27 (ArC), 145.26 (ArC), 144.69 (ArC), 139.48 (ArC), 134.82 (ArC), 129.48 (ArC), 127.42 (ArC), 125.50 (ArC), 121.50 (ArC), 120.13 (ArC), 118.09 (ArC), 117.08 (ArC), 115.14 (ArC), 112.76 (ArC), 112.59 (ArC), 109.63 (ArC); LC/MS Calculated C₁₈H₁₃N₃O₂ [M+1]⁺: m/z 303.32; observed, 304.10.

Table S1: Physical properties of imidazopyridine amides **6a-r**

Molecule Id	R	Molecular weight	R_f	% Yield
6a	-H	313.36	0.40	80.29
6b	-4F	331.35	0.41	85.23
6c	-4Cl	347.80	0.44	89.56
6d	-4Br	392.26	0.47	88.64
6e	-4CF ₃	381.36	0.39	87.03
6f	-4CH ₃	327.39	0.30	78.11
6g	-4OCH ₃	343.39	0.29	72.36
6h	-4NO ₂	358.36	0.38	76.25
6i	-3F	331.35	0.39	82.56
6j	-3Cl	347.80	0.45	88.44
6k	-3Br	392.26	0.36	78.66
6l	-3CF ₃	381.36	0.34	85.77
6m	-3CH ₃	327.39	0.38	75.96
6n	-3OCH ₃	343.39	0.31	71.20
6o	-3NO ₂	358.36	0.33	86.55
6p	-2F	331.35	0.42	84.65
6q	1-Napthyl	363.42	0.45	77.59
6r	2-Furyl	303.32	0.32	69.98