**Supplementary Tables**

**Table S1**. Characteristic constants of fluorinated pyridine Schiff bases **F1** and **F2**.

|  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- |
| **Schiff Base** | **Molecular Weight (g mol-1)** | **Yield (%)** | **Melting point (ºC)** | **Solid Color** | **Thin Layer Chromatography Rf**  |
| **F1** | 249.0 | 78 | 182.5 – 183.7 | Yellow | 0.31 |
| **F2** | 231.0 | 82 | 163.4 – 164.8 | Yellow | 0.40 |

**Table S2.** UV-Vis absorption spectra of **F1** and **F2**.

|  |  |  |  |
| --- | --- | --- | --- |
| **Compound** | **Dichloromethane** | **Acetonitrile** | **DMSO** |
|  | **λ exp\*** | **ε**  | **λ exp** | **ε**  | **λ exp** | **ε** |
| **F1** | 366260 (sh 278) | 6.2911.59 | 362 | 7.76 | 374262 | 7.7014.96 |
| **F2** | 364274 | 9.4610.18 | 360 | 9.30 | 372262 | 9.9015.80 |

\* λ in nm

ε (103 mol-1 dm3 cm-1)

**Table S3.** Optimized geometrical parameters (All the distances are in Å).

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
| **Compound** | **d(N-Cpy)** | **d(C-N)****azomethine group** | **d(N-H)** | **d(O-H)****phenolic ring** |
| **F1** | 1.367 | 1.284 | 1.772 | 0.989 |
| **F2** | 1.368 | 1.285 | 1.773 | 0.988 |

**Table S4.** Frequencies of the principal groups calculated for **F1** and **F2**.

|  |  |  |  |
| --- | --- | --- | --- |
| **Compound** | **νOH (cm-1)** | **νNH (cm-1)** | **νN=C (cm-1)** |
| **F1** | 3252 | νas: 3702νs: 3595 | 1626 |
| **F2** | 3252 | νas: 3701νs: 3593 | 1592 |

**Table S5.** Most important transition energies calculated for **F1** and **F2**

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
| **Compound** | **Solvent** | **λ(nm)** | **f** | **Assignment** |
| **F1** | CH2Cl2 | 216 | 0.32 | HOMO-1 → LUMO+4 (n→π\*)HOMO-2 → LUMO+1 (n→π\*) |
| 288 | 0.29 | HOMO-3 → LUMO (n→π\*)HOMO-2 → LUMO (n→π\*) |
| 390 | 0.47 | HOMO → LUMO (π→π\*) |
| ACN | 214 | 0.28 | HOMO-1 → LUMO+4 (n→π\*)HOMO-2 → LUMO+1 (n→π\*) |
| 289 | 0.26 | HOMO-3 → LUMO (n→π\*)HOMO-2 → LUMO (n→π\*) |
| 385 | 0.30 | HOMO → LUMO (π→π\*) |
| DMSO | 220 | 0.38 | HOMO-1 → LUMO+4 (n→π\*)HOMO-2 → LUMO+1 (n→π\*) |
| 284 | 0.31 | HOMO-3 → LUMO (n→π\*)HOMO-2 → LUMO (n→π\*) |
| 390 | 0.30 | HOMO → LUMO (π→π\*) |
| Gas phase | 213 | 0.28 | HOMO-1 → LUMO+3 (n→π\*)HOMO-2 → LUMO+1 (n→π\*) |
| 277 | 0.24 | HOMO-3 → LUMO (n→π\*)HOMO-2 → LUMO (n→π\*) |
| 394 | 0.33 | HOMO → LUMO (π→π\*) |
| **F2** | CH2Cl2 | 220 | 0.30 | HOMO-2 → LUMO+3 (n→π\*)HOMO-2 → LUMO+1 (n→π\*) |
| 290 | 0.32 | HOMO-3 → LUMO (n→π\*)HOMO-2 → LUMO (n→π\*) |
| 380 | 0.56 | HOMO → LUMO (π→π\*) |
| ACN | 215 | 0.40 | HOMO-1 → LUMO+3 (n→π\*)HOMO-2 → LUMO+1 (n→π\*) |
| 283 | 0.30 | HOMO-3 → LUMO (n→π\*)HOMO-2 → LUMO (n→π\*) |
| 387 | 0.56 | HOMO → LUMO (π→π\*) |
| DMSO | 218 | 0.40 | HOMO-1 → LUMO+3 (n→π\*)HOMO-2 → LUMO+1 (n→π\*) |
| 285 | 0.32 | HOMO-3 → LUMO (n→π\*)HOMO-2 → LUMO (n→π\*) |
| 382 | 0.56 | HOMO → LUMO (π→π\*) |
| Gas phase | 216 | 0.22 | HOMO-1 → LUMO+3 (n→π\*)HOMO-2 → LUMO+1 (n→π\*) |
| 284 | 0.27 | HOMO-3 → LUMO (n→π\*)HOMO-2 → LUMO (n→π\*) |
| 389 | 0.47 | HOMO → LUMO (π→π\*) |

**Table S6.** Scan rate study results

|  |  |  |
| --- | --- | --- |
| Compound | **F1** | **F2** |
| Process | Eox | Ered | Eox  | Ered |
| V vs SCE | +1.07 | -0.89 | +0.92 | -1.31 |
| Scan-rate vs current-density peak | m | 6.0×107 | -3.0×107 | 5.0×106 | -2.0×106 |
| n | -96.3 | -131.2 | 1.4 | -7.0 |
| R2 | 0.914 | 0.968 | 0.990 | 0.882 |
| (Scan-rate)½ vs current-density peak | m | 3.0×106 | -2.0×106 | 9.0×107 | -5.0×107 |
| n | -1.0 | -2.6 | -56.0 | -201.5 |
| R2 | 0.973 | 0.996 | 0.997 | 0.786 |
| Diffusion control? | yes | yes | no | yes |

**Table S7.** Minimal inhibition concentration (μg/mL) of tested compounds alone or included in **βCD**

|  |  |
| --- | --- |
|  | **Concentration (μg/mL) ± SE** |
| **Compound** | ***Salmonella enterica*****(18 h)** | ***Staphylococcus aureus*****(18 h)** |
| **F1** | - | - |
| **F2** | - | - |
| **βF1** | NE | 200.0 ± 0.0 |
| **βF2** | NE | NE |
| **βCD** | NE | NE |

SE: Standard error

- : Undistinguishable from DMSO alone (i.e. no effect)

NE: No effect