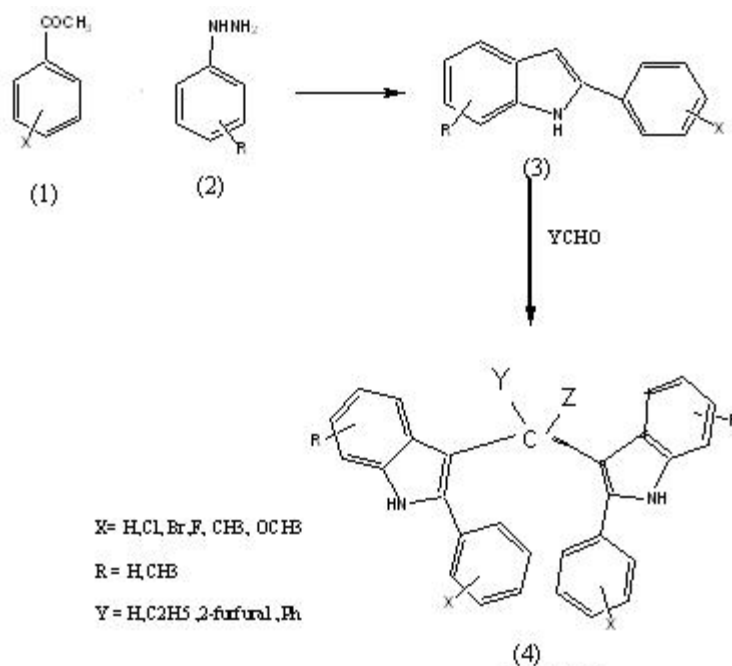


### 3. RESULT AND DISCUSSION

Title compounds (4a-j) were obtained by refluxing indole and appropriate aldehyde or ketone in 2:1 ratio by scheme-1 as shown in Figure 1.



**Fig. 1:** Scheme-1 to prepare title compounds.

The IR spectra of title compounds (4a-j) showed strong >N-H absorption band from 3417-3400  $\text{cm}^{-1}$  which was present earlier in compound (3) and aromatic C-H str. absorption peak appears at 3060-3051  $\text{cm}^{-1}$ . Aliphatic C-H stretching Vibration is observed at 2839  $\text{cm}^{-1}$  and aromatic C=C absorption band is observed at 1604  $\text{cm}^{-1}$ . In their PMR all the diindolylmethanes exhibited disappearance of singlet from  $\delta$  6.4ppm assigned to methine proton (=C-H, at C-3) of the starting arylindole (3). Compound 4(a) showed additional resonance signal as double doublet from  $\delta$  4.02 - 4.52 ppm due to  $\text{CH}_2$  protons indicating that they are nonequivalent protons. Aromatic proton (Ar-H) appears as a complex multiplet in the region of  $\delta$  6.8-7.76 ppm. Additional characteristic peaks corresponding to the formation of diindolylmethane were recorded in Table 2. Mass spectral data were also satisfactory and corresponded to their molecular weights as shown in Table 2.

**Table 2:** Spectral data of 2- phenyl indole (3) and diindolyl methan 4a,4c,4e,4f

| Com pd. No. | X                 | Y                             | 1 R (KBr) $\nu_{\max}$ $\text{cm}^{-1}$   | $^1\text{H NMR}$ $\delta$ ppm   | Mass m/z       |
|-------------|-------------------|-------------------------------|---|---|----------------|
| 3           | H                 | -                             | 3450–3350 (br, N,H), 1610 (C = C)   | 6.4(s,CH,1H),6.8- 8.0 (m,NH+ArH,10H)                                  | 193 (M+)       |
| 4a          | F                 | H                             | 3408(NHstr.), 3051 (aromatic C-H str.),2839 (aliphatic C-H str.) 1683(aromatic C=C str.), 1603(C = C str.Attached to N)                 | 4.02-4.52(dd,CH <sub>2</sub> ,2H),6.80-7.47(m,ArH,16 H),7.76(s,NH,2H) | 434 (M+)       |
| 4c          | Br                | Furfural                      | 3412 (NH str), 3026 (aromatic C-H str.),2840 (aliphatic C-H str.)1685 (aromatic C=C str.), 1604(C=C str. Attached to N) 750 (C-Br str.) | 6.26(s,CH,1H),6.78-7.75(m,ArH,21H), 8.26(s,NH,2H)                     | 606 (M+2) peak |
| 4e          | F                 | C <sub>6</sub> H <sub>5</sub> | 3408 (NH str.),3030 (aromatic C-H str.), 2832(aliphatic C-H str), 1690(aromatic C=C str), 1602(C=Cstr attached to N)                    | 6.61(s,CH,1H),6.88-7.65(m,ArH,21H), 8.26(s,NH,2H)                     | 511 (M+)       |
| 4f          | OC H <sub>3</sub> | C <sub>6</sub> H <sub>5</sub> | 3415(NHstr.),3049 (aromatic C-H str.),2839 (aliphatic C-H str), 1683(aromatic C=C str.), 1602(C=C str. Attached to N)                   | 6.04(s,CH,1H),6.86-7.36(m,ArH,21H), 7.92(s,NH,2H)                     | 534 (M+)       |

### 3.1. Antibacterial Activities

Compounds (4a-j) were screened against gram negative bacteria *Escherichia coli* and gram positive bacteria *klebsiella* at different concentrations by disk diffusion method [15] as tabulated in Table 3. Compound (4c,4f,4h) showed activity greater than standard activity against *E.Coli* and *klebsiella* at 1000 ppm concentration.

**Table 3:** Antibacterial activity of 2-phenylindole (3) and diindolylalkane (4a-j)

| Compd No. | Mean value of area of inhibition in mm (1000ppm) |            | Mean value of area of inhibition in mm (800ppm) |            | Mean value of area of inhibition in mm (400ppm) |            | Mean value of area of inhibition in mm (200ppm) |            |
|-----------|--|------------|---|------------|---|------------|---|------------|
|           | E.Co li  | klebsiella | E.Co li   | klebsiella | E.Co li   | klebsiella | E.Col i   | klebsiella |
| Amicacn   | 12   | 13         | 10  | 11         | 08  | 10         | 07  | 06         |
| 3         | 08   | 10         | 08  | 10         | 06  | 07         |   | 03         |
| 4a        | 06   | 12         | 04  | 08         | 01  | 02         | nil   | 04         |
| 4b        | 10   | 14         | 06  | 10         | 04  | 04         |   | 02         |
| 4c        | 16   | 12         | 12  | 10         | 08  | 06         | 04  | 03         |
| 4d        | 09   | 10         | 10  | 08         | -   | 06         | -   | 01         |
| 4e        | 07   | 08         | 10  | 08         |   | 04         | -   | 02         |
| 4f        | 14   | 12         | 10  | 09         | 18  | 16         | -   | 03         |
| 4g        | 08   | 07         | nil   | 02         | 06  | nil        | -   | nil        |
| 4h        | 27   | 16         | 13  | 07         | 14  | 04         | 08  | 04         |
| 4i        | 11   | 12         | 10  | 14         | 06  | -          | 04  | -          |
| 4j        | 06   | 06         | 04  | 02         | 03  | nil        | 02  | nil        |

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