

h. IR (KBr) Spectral data (ν_{max} in cm^{-1})

VIIa: 3206 (NH), 1685 (C=O), 1620 (C=N)

VIIb: 3195 (NH), 1690 (C=O), 1622 (C=N)

VIIc: 3230 (NH), 1685 (C=O), 1625 (C=N)

VIId: 3215 (NH), 1695 (C=O), 1624 (C=N)

VIIe: 3230 (NH), 1675 (C=O), 1630 (C=N)

VIIf: 3210 (NH), 1685 (C=O), 1627 (C=N)

VIIg: 3180 (NH), 1695 (C=O), 1627 (C=N)

VIIh: 3195 (NH), 1700 (C=O), 1629 (C=N)

VIIi: 3245 (NH), 1705 (C=O), 1630 (C=N)

VIIj: 3240 (NH), 1685 (C=O), 1630 (C=N)

i. 1H NMR (DMSO- d_6) Spectral data (δ in ppm)VIIa: 2.22 (s, 3H CH₃), 2.40 (s, 3H, CH₃), 2.46 (s, 3H, COCH₃), 5.26 (s, 2H, OCH₂), 6.8 (s, H, Ar-NH), 7.1-7.3 (m, 9H, Ar-H), 7.4 (d, 2H, C₆H₄), 7.7 (d, 2H, C₆H₄).VIIg: 2.20 (s, 3H CH₃), 2.26 (s, 3H, CH₃), 2.39 (s, 3H, CH₃), 2.44 (s, 3H, COCH₃), 5.24 (s, 2H, OCH₂), 6.8 (s, H, Ar - NH), 7.1-7.3 (m, 9H, Ar-H), 7.4 (d, 2H, C₆H₄), 7.7 (d, 2H, C₆H₄).VIIh: 2.19 (s, 3H, CH₃), 2.37 (s, 3H, COCH₃), 2.24 (s, 2H, CH₃), 4.92 (s, 2H, OCH₂), 6.8 (s, H, Ar-NH), 7.1-7.3 (m, 9H, Ar-H), 7.4 (d, 2H, C₆H₄), 7.7 (d, 2H, C₆H₄).VIIi: 2.18 (s, 3H CH₃), 2.39 (s, 3H, CH₃), 2.43 (s, 3H, COCH₃), 3.89 (s, 3H, OCH₃), 5.24 (s, 2H, OCH₂), 6.8 (s, H, Ar-NH), 7.1-7.3 (m, 9H, Ar-H), 7.4 (d, 2H, C₆H₄), 7.7 (d, 2H, C₆H₄).VIIj: 2.16 (s, 3H CH₃), 2.35 (s, 3H, CH₃), 2.40 (s, 3H, COCH₃), 5.22 (s, 2H, OCH₂), 6.8 (s, H, Ar - NH), 7.1-7.3 (m, 9H, Ar-H), 7.4 (d, 2H, C₆H₄), 7.7 (d, 2H, C₆H₄).**j. Elemental analysis data of compounds VIII (M.P., Yield, Molecular formula, Element: Found %, (Calc %))**VIII: 150 °C, 65 %, C₁₉H₂₀N₆O₃S, C:55.33(55.19), H:4.85(4.68), N:20.38(20.21), O:16.50(16.35), S:7.76(7.62).**k. IR (KBr) Spectral data (ν_{max} in cm^{-1})**

VIII: 3126 (oxadiazole NH), 3180 (NH), 1603 (C=N), 1670 (C=O), 1134 (C=S).

l. 1H NMR (DMSO- d_6) Spectral data (δ in ppm)VIII: 2.3 (s, 3H CH₃), 5.45 (s, 2H, OCH₂), 6.8 (s, H, Ar - NH), 7.1-7.3 (m, 5H, Ar - H), 7.4 (d, 2H, C₆H₄), 7.7 (d, 2H, C₆H₄), 14.7 (s, H, thiol-thione tautomeric proton NH)**m. Elemental analysis data of compounds IX (R₁=H, R₂, M.W., M.P., Yield, Molecular formula, Element: Found %, (Calc %))**IXa: *p*-tolyl, 240°C, 75 %, C₂₇H₂₅N₇O₃S, C:61.48(61.32), H:4.47(4.30), N:18.59(18.45), O:9.10(8.93), S:6.07(5.90).IXb: *p*-anisyl, 245°C, 77 %, C₂₇H₂₅N₇O₄S, C:59.66(59.48), H:4.60(4.56), N:18.04(17.85), O:11.78(11.62), S:5.89(5.69).IXc: *p*-fluorophenyl, 235°C, 78 %, C₂₆H₂₂N₇O₃SF, C:58.76(58.65), H:4.14(3.98), N:18.45(18.33), O:9.04(8.86), S:6.02(5.80), F:3.55 (3.38).IXd: *p*-chlorophenyl, 250°C, 73 %, C₂₆H₂₂N₇O₃SCl, C:56.98(56.75), H:4.01(3.83), N:17.88(17.69), O:8.76(8.61), S:5.84(5.69), Cl:6.48(6.31).IXe: *p*-bromophenyl, 230°C, 80 %, C₂₆H₂₃N₇O₃Br, C:52.71(52.58), H:3.71(3.59), N:16.55(16.38), O:8.10(7.92), S:5.40(5.23), Br:13.49(13.30).IXf: *p*-nitrophenyl, 255°C, 83 %, C₂₆H₂₂N₈O₅S, C:55.91(55.74), H:3.94(3.78), N:20.07(19.87), O:14.33(14.21), S:5.73(5.57).IXg: diethyl, 260°C, 72 %, C₂₈H₃₇N₈O₃S, C:59.46(59.28), H:6.54(6.39), N:19.82(19.70), O:8.49(8.32), S:5.66(5.44).IXh: diphenyl, 266 °C, 70 %, C₃₂H₂₇N₇O₃S, C:65.29(65.18), H:4.52(4.62), N:17.01(16.63), O:8.03(8.14), S:5.27(5.44).IXi: morpholinyl, 270°C, 70 %, C₂₀H₁₈N₇O₄S, C:53.09(52.89), H:3.98(3.76), N:21.68(21.52), O:14.15(13.98), S:7.07(6.84).IXj: piperazinyl, 272°C, 72 %, C₂₂H₂₀N₇O₃S, C:57.14(56.98), H:4.32(4.17), N:21.21(21.07), O:10.38(10.23), S:6.92(6.76).