Reactions with pyridine. In the pyridine molecule there is only one nitrogen atom situated in the plane parallel to the CR'R"R"" plane and perpendicular to the ring (Fig. 2). Two "ortho"-hydrogens do not reach this plane (by $\sim 0.2\text{Å}$), and two "ortho"-carbons - by 0.7Å. A compression originates, in the reaction with methyl iodide, between nitrogen atom and three hydrogen atoms (1.85Å), between one H in CH₃I and an H atom in C₅H₅N (1.7Å instead of 2.4Å) and also between C atoms in C₃H₅N and H atoms in CH₃I. The sum of these compressions is equal to 6.7Å³, i.e. 4 cc/mol.; $\triangle v_{\text{mol.}}^{\#} = 9$ cc/mol.

In the reaction with ethyl iodide, one compression N...H is changed into N...CH₃ (2.15Å instead of 3.6Å) and there arise interactions with "ortho"- carbons; $\triangle \sqrt[4]{mol}$. == 14 cc/mol.

The value $\Delta \vec{v}_{\text{mol.}}$ does not change with further increase of the number of carbon atoms in the normal little chain of alkyl halide.

In the reaction with <u>iso</u>propyl iodide, there are two N...CH₃ compressions and the character of the interaction with "ortho"-carbons is somewhat changed, t mol. =-19 cc/mol.

Reactions with trimethylamine. Consider the reaction of trimethylamine with isopropyl iodide. The most preferable mutual disposition of the CR3 plane and N(CH3)3 pyramide is represented by the scheme in Fig. 3a (along the C-I bond).

In the activated complex originate the following compressions: compression N...H (1.85Å), two $N...CH_3$ compressions (2.6Å) and four $CH_3...CH_3$ compressions (2.7Å). The calculated value of $\Delta_{\mathcal{D}}^{+}$ mol. is equal to -24 cc/mol., i.e. larger than in the reaction of isopropy! iodide with pyridine. This value does not

^{**} Including the compression in the formation of a C-N bond (see above).